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Certifications

WBENC: 237019

HUB:

1752439743100-86536

DBE: VN 20657

NCTRCA WFWB38444Y0909

NELAP Certifications

Lubbock: T104704219-08-TX El Paso:

T104704221-08-TX

Midland: T104704392-08-TX

LELAP-02003 Kansas E-10317 LELAP-02002

Analytical and Quality Control Report

Brad Davis Zia Engineering & Environmental 755 S. Telshor Blvd. Suite F-201 Las Cruces, NM, 88011

Report Date: September 18, 2009

Work Order:

9090211

Project Name: HELSTF Diesel Spill Groundwater

Enclosed are the Analytical Report and Quality Control Report for the following sample(s) submitted to TraceAnalysis,

Inc.

			Date	Time	Date
\mathbf{Sample}	Description	Matrix	Taken	Taken	Received
208954	HLSF-0154-DRW-112-0809	water	2009-08-31	10:57	2009-08-31

Comment(s)

These results represent only the samples received in the laboratory. The Quality Control Report is generated on a batch basis. All information contained in this report is for the analytical batch(es) in which your sample(s) were analyzed.

This report consists of a total of 80 pages and shall not be reproduced except in its entirety, without written approval of TraceAnalysis, Inc.

Notes:

For inorganic analyses, the term MQL should actually read PQL.

Standard Flags

- ${f U}\,$ Not detected. The analyte is not detected above the SDL.
- ${f J}$ Estimated. The analyte is positively identified and the value is approximated between the SDL and MQL.
- B The sample contains less than ten times the concentration found in the method blank.
- JB The analyte is positively identified and the value is approximated between the SDL and MQL.

The sample contains less than ten times the concentration found in the method blank.

The result should be considered non-detect to the SDL.

Dr. Blair Leftwich, Director
Dr. Michael Abel, Project Manager

Case Narrative

Samples for project HELSTF Diesel Spill Groundwater were received by TraceAnalysis, Inc. on 2009-08-31 and assigned to work order 9090211. Samples for work order 9090211 were received intact without headspace and at a temperature of 12.0 deg. C, just sampled, on ice.

Samples were analyzed for the following tests using their respective methods.

		Prep	Prep	QC	Analysis
Test	Method	Batch	Date	Batch	Date
Ag, Total	S 6010B	53951	2009-09-03 at 08:16	63221	2009-09-03 at 10:42
${ m Alkalinity}$	SM 2320B	54231	2009-09-10 at $11:00$	63527	2009-09-10 at 11:00
Al, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
${ m Ammonia}$	SM 4500-NH3 B,C	54092	2009-09-05 at $16:00$	63370	2009-09-05 at $17:00$
As, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Ba, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Be, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Bromide (IC)	E 300.0	54363	2009-09-16 at $13:48$	63674	2009-09-16 at $13:48$
Ca, Total	S 6010B	53951	2009-09-03 at $08:16$	63355	2009-09-09 at $08:18$
Cd, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Chloride (IC)	E 300.0	54356	2009-09-01 at $23:46$	63668	2009-09-01 at $23:46$
Chromium, Hexavalent	$\mathrm{SM}\ 3500\text{-Cr}\ \mathrm{B}$	54055	2009-09-01 at $10:12$	63327	2009-09-01 at $10:12$
Co, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Cr, Dissolved	S 6010B	54153	2009-09-11 at $08:25$	63461	2009-09-11 at 11:54
Cr, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Cu, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Explosives (8330)	S 8330-C18	54137	2009-09-04 at 15:00	63425	2009-09-10 at 15:30
Fe, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Fluoride (IC)	E 300.0	54363	2009-09-16 at $13:48$	63674	2009-09-16 at $13:48$
Hg, Total	S 7470A	53960	2009-09-03 at $12:02$	63227	2009-09-03 at $13:18$
K, Total	S 6010B	53951	2009-09-03 at $08:16$	63355	2009-09-09 at $08:18$
Mg, Total	S 6010B	53951	2009-09-03 at $08:16$	63355	2009-09-09 at $08:18$
Mn, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Mo, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Na, Total	S 6010B	53951	2009-09-03 at $08:16$	63355	2009-09-09 at $08:18$
Ni, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Nitrate and Nitrite as N	SM 4500-NO3 E	54188	2009-09-11 at $10:24$	63485	2009-09-11 at $16:25$
O/G	$\to 1664$	54391	2009-09-17 at 11:00	63699	2009-09-17 at 14:30
Pb, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
pН	SM 4500-H+	54063	2009-09-01 at $11:30$	63339	2009-09-01 at $11:30$
P, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Sb, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Semivolatiles	S 8270C	53996	2009-09-02 at $15:00$	63261	2009-09-04 at $10:45$
Se, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
SO4 (IC)	E 300.0	54356	2009-09-01 at 23:46	63668	2009-09-01 at $23:46$
TDS	SM 2540C	54173	2009-09-03 at $14:20$	63473	2009-09-03 at $14:20$
TKN	$\to 351.3$	54150	2009-09-09 at $11:15$	63441	2009-09-09 at $15:00$
Tl, Total	S 6010B	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
TOC	SM 5310C	54065	2009-09-08 at $13:26$	63340	2009-09-08 at $13:29$
Total Cyanide	SM 4500-CN C, $\rm E$	54106	2009-09-07 at $16:00$	63391	2009-09-07 at 17:45
TPH DRO	Mod. 8015B	54035	2009-09-04 at 15:00	63307	2009-09-07 at 18:00

		Prep	Prep	QC	${ m Analysis}$
Test	Method	Batch	Date	Batch	Date
TPH GRO	S~8015B	53976	2009-09-03 at 14:48	63239	2009-09-03 at $14:48$
V, Total	S_{010B}	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$
Zn, Total	S_{6010B}	53951	2009-09-03 at $08:16$	63221	2009-09-03 at $10:42$

Results for these samples are reported on a wet weight basis unless data package indicates otherwise.

A matrix spike (MS) and matrix spike duplicate (MSD) sample is chosen at random from each preparation batch. The MS and MSD will indicate if a site specific matrix problem is occurring, however, it may not pertain to the samples for work order 9090211 since the sample was chosen at random. Therefore, the validity of the analytical data reported has been determined by the laboratory control sample (LCS) and the method blank (MB). These quality control measures are performed with each preparation batch to ensure data integrity.

All other exceptions associated with this report have been footnoted on the appropriate analytical page to assist in general data comprehension. Please contact the laboratory directly if there are any questions regarding this project.

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HELSTF Diesel Spill Groundwater

Analytical Report

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analytical Method: Analysis: Ag, Total S 6010B Prep Method: S 3010A RRQC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: Prep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQLMethod Based Based Blank MQLMDL SDLParameter Result Result Result Dilution (Unadjusted) (Unadjusted) Flag Units Total Silver 0.00111 < 0.00111 < 0.00500< 0.00111mg/L0.00111 0.005

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Al, Total S 3010A Analysis: Analytical Method: S 6010B Prep Method: QC Batch: Date Analyzed: 2009-09-03 Analyzed By: 63221 RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQLMethod Based Based Blank MQLMDLParameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Total Aluminum 0.2370.237< 0.00301 mg/L0.003010.050.00301

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: El Paso

Analysis: Alkalinity Analytical Method: SM 2320BPrep Method: N/AQC Batch: 63527 Date Analyzed: 2009-09-10 Analyzed By: JGPrep Batch: Sample Preparation: Prepared By: 54231JG

SDLMQL Method Based Based Blank MQLMDL Flag Parameter Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Hydroxide Alkalinity < 1.00 <1.00 < 1.00 mg/L as CaCo3 1.00 1 UCarbonate Alkalinity < 1.00 < 1.00 < 1.00mg/L as CaCo3 1 1.00 1 1 Bicarbonate Alkalinity 264 264 < 4.00 mg/L as CaCo3 4.00 4 4 1 264 mg/L as CaCo3Total Alkalinity 264 < 4.001 4.00 4 4

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Analytical Method: Prep Method: N/AAmmonia SM 4500-NH3 B,C QC Batch: 63370 Date Analyzed: Analyzed By: AH2009-09-05 2009-09-05 Prep Batch: 54092 Sample Preparation: Prepared By: AH

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HELSTF Diesel Spill Groundwater

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Ammonia-N	J	0.560	<1.00	< 0.353	mg/L	1	0.353	1	0.353

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: As, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 2009-09-03 63221 Date Analyzed: Analyzed By: RRPrep Batch: Sample Preparation: 2009-09-03 Prepared By: KV53951

SDLMQLMethod Based Based Blank MQLMDL Parameter Result Result Units Dilution SDLFlag Result (Unadjusted) (Unadjusted) Total Arsenic 0.01200.0120< 0.00448 mg/L0.00448 0.01 0.00448 1

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Ba, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: Sample Preparation: 2009-09-03 Prepared By: KV53951

SDLMQLMethod MDL Based Based Blank MQLParameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Total Barium 0.01600.0160< 0.00105 0.001050.005 0.00105 mg/L1

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Be, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: Date Analyzed: 2009-09-03 Analyzed By: RR63221 Prep Batch: 53951Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQLMethod MQLMDL Based Based Blank Parameter Flag Result Result Result Dilution SDL(Unadjusted) Units (Unadjusted) Total Beryllium < 0.000450 < 0.00200 < 0.000450 mg/L0.0004500.002 0.00045

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: El Paso

Analysis: Bromide (IC) Analytical Method: E 300.0 Prep Method: N/A

Report Date: September 18, 2009 Work Order: 9090211 Page Number: 7 of 80

HELSTF Diesel Spill Groundwater

QC Batch: Prep Batch:	63674 54363			Date Analyzed: Sample Preparation:		2009-09-16 2009-09-16		Analyzed By: JR Prepared By: JR		
		SDL	MQL	Method						
		Based	Based	Blank				MQL	MDL	
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)	
Bromide		3.33	3.33	< 0.197	$_{ m mg/L}$	5	0.197	0.27	0.0394	

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Ca, Total Analysis: Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQLMethod Based Based Blank MQLMDL Parameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Total Calcium 447 447< 1.17mg/L10 1.17 0.117 1

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Cd, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: Analyzed By: 63221 Date Analyzed: 2009-09-03 RRPrep Batch: 53951Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQLMethod Based Based Blank MQLMDL Parameter Flag Result Result ResultUnits Dilution SDL(Unadjusted) (Unadjusted) Total Cadmium < 0.000303 < 0.00200 < 0.000303 0.0003030.0020.000303mg/L

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: El Paso

Analysis: Chloride (IC) Analytical Method: E 300.0 Prep Method: N/AQC Batch: 63668 Date Analyzed: 2009-09-01 Analyzed By: JRPrep Batch: Prepared By: JR54356Sample Preparation: 2009-09-01

SDLMQLMethod Based Based Blank MQL MDLSDLParameter Flag Result Result Result Units Dilution (Unadjusted) (Unadjusted) Chloride 4920 4920 <320 mg/L500 320 1.22 0.6404

Sample: 208954 - HLSF-0154-DRW-112-0809

Report Date: September 18, 2009 Work Order: 9090211 Page Number: 8 of 80

HELSTF Diesel Spill Groundwater

Laboratory: El Paso

Analysis: Analytical Method: Chromium, Hexavalent SM 3500-Cr B Prep Method: N/AQC Batch: Date Analyzed: 2009-09-01 Analyzed By: MD63327 Prep Batch: 54055Sample Preparation: 2009-09-01 Prepared By: JR

> SDLMQLMethod

Based Based Blank MQL MDL Parameter Flag Result Result ${\bf Result}$ Units Dilution SDL (Unadjusted) (Unadjusted) Hexavalent Chromium 0.01040.0104< 0.00594mg/L0.005940.01 0.005941

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Co, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQL Method Based Based Blank

MQLMDLParameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Total Cobalt < 0.000822 < 0.00200 < 0.000822 mg/L0.0008220.0020.000822

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Analytical Method: Prep Method: S 3005A Cr, Dissolved S 6010B QC Batch: 63461 Date Analyzed: 2009-09-11 Analyzed By: RRPrep Batch: 54153Sample Preparation: 2009-09-11 Prepared By: KV

SDLMQLMethod Based Based Blank

MQLMDL SDLResult Result Parameter Flag Result Dilution (Unadjusted) (Unadjusted) Units Dissolved Chromium 0.01900.000583 0.0190< 0.000583 mg/L0.0005830.001

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Cr, Total Analytical Method: Prep Method: S 3010A S 6010B QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDL \mathbf{MQL} Method

		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Chromium		0.572	0.572	< 0.000583	$\mathrm{mg/L}$	1	0.000583	0.005	0.000583

HELSTF Diesel Spill Groundwater

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Cu, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR Prep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDL MQL Method

		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Copper		0.0230	0.0230	< 0.000843	mg/L	1	0.000843	0.005	0.000843

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Explosives (8330) Analytical Method: S 8330-C18 Prep Method: S 3535A QC Batch: 63425Date Analyzed: 2009-09-10 Analyzed By: DSPrep Batch: 54137Sample Preparation: 2009-09-04 Prepared By: DS

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	$\operatorname{Dilution}$	SDL	(Unadjusted)	(Unadjusted)
HMX	1 U	< 0.246	<1.00	< 0.246	$\mu \mathrm{g/L}$	2	0.246	0.5	0.123
RDX	U	< 0.596	< 1.00	< 0.596	$\mu { m g/L}$	2	0.596	0.5	0.298
1,3,5-Trinitrobenzene	U	< 0.678	< 1.00	< 0.678	$\mu { m g/L}$	2	0.678	0.5	0.339
1,3-Dinitrobenzene	U	< 0.778	< 1.00	< 0.778	$\mu { m g/L}$	2	0.778	0.5	0.389
Nitrobenzene	U	< 0.758	< 1.00	< 0.758	$\mu \mathrm{g/L}$	2	0.758	0.5	0.379
Tetryl	U	< 0.826	< 1.00	< 0.826	$\mu { m g/L}$	2	0.826	0.5	0.413
TNT	U	< 0.928	< 1.00	< 0.928	$\mu \mathrm{g/L}$	2	0.928	0.5	0.464
4-Amino-DNT	U	< 0.638	< 1.00	< 0.638	$\mu \mathrm{g/L}$	2	0.638	0.5	0.319
2-Amino-DNT	U	< 0.782	< 1.00	< 0.782	$\mu { m g}/{ m L}$	2	0.782	0.5	0.391
2,6-DNT	U	< 0.646	< 1.00	< 0.646	$\mu \mathrm{g/L}$	2	0.646	0.5	0.323
2,4-DNT	U	< 0.732	< 1.00	< 0.732	$\mu \mathrm{g}/\mathrm{L}$	2	0.732	0.5	0.366
2-NT	U	< 0.758	< 1.00	< 0.758	$\mu \mathrm{g/L}$	2	0.758	0.5	0.379
4-NT	U	< 0.796	< 1.00	< 0.796	$\mu \mathrm{g}/\mathrm{L}$	2	0.796	0.5	0.398
3-NT	U	< 0.692	< 1.00	< 0.692	$\mu \mathrm{g}/\mathrm{L}$	2	0.692	0.5	0.346

					Spike	$\operatorname{Percent}$	$\operatorname{Recovery}$
$\mathbf{Surrogate}$	Flag	Result	Units	Dilution	${f Amount}$	Recovery	Limits
1,2-Dinitrobenzene		1.79	$\mu \mathrm{g/L}$	2	2.50	72	19.8 - 160

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Fe, Total Analytical Method: Prep Method: S 3010A S 6010B QC Batch: 63221Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

¹Sample ran at a dilution due to matrix difficulties.

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HELSTF Diesel Spill Groundwater

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	$\operatorname{Dilution}$	SDL	(Unadjusted)	(Unadjusted)
Total Iron		3.50	3.50	< 0.000872	mg/L	1	0.000872	0.01	0.000872

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: El Paso

Analysis: Fluoride (IC) Analytical Method: E 300.0 Prep Method: N/AQC Batch: 63674 Date Analyzed: 2009-09-16 Analyzed By: JRPrep Batch: 54363 Sample Preparation: Prepared By: JR2009-09-16 SDLMQLMethod Based Based Blank MQLMDL Parameter Result Units Dilution SDLFlag Result Result (Unadjusted) (Unadjusted) Fluoride 2.572.57< 0.217mg/L5 0.2170.17 0.0434

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Hg, Total Analytical Method: S 7470A Prep Method: N/AQC Batch: TP63227 Date Analyzed: 2009-09-03 Analyzed By: Prep Batch: Sample Preparation: 2009-09-03 Prepared By: TP 53960

SDLMQL Method MQLMDL Based Based Blank Parameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Total Mercury < 0.0000329 < 0.000200 < 0.0000329 0.00003290.0002 3.29e-05mg/L

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: K, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RRPrep Batch: 53951Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQLMethod MQLMDL Based Based Blank SDL Parameter Flag Result Result Result Units Dilution (Unadjusted) (Unadjusted) Total Potassium 216 216 < 1.72mg/L10 1.72 0.172

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Mg, Total Analytical Method: S 6010B Prep Method: S 3010A

Report Date: September 18, 2009 Work Order: 9090211 Page Number: 11 of 80

HELSTE	Diesel	Spill	Groundwater
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QC Batch: Prep Batch:				Date Analyzed: Sample Preparation:		9-09-09 9-09-03		Analyzed E Prepared B	•	
			SDL	MQL	Method					
			Based	Based	Blank				MQL	MDL
Parameter		Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Magnes	sium		1710	1710	<16.0	$\mathrm{mg/L}$	100	16.0	1	0.16

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory:	$\operatorname{Lubbock}$
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Analysis:	Mn, Total	Analytical Method:	S 6010B	Prep Method:	S 3010A
QC Batch:	63221	Date Analyzed:	2009-09-03	Analyzed By:	RR
Prep Batch:	53951	Sample Preparation:	2009-09-03	Prepared By:	KV

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Manganese		0.0260	0.0260	< 0.000305	mg/L	1	0.000305	0.0025	0.000305

Sample: 208954 - HLSF-0154-DRW-112-0809

La	boratory:	Lul	b]	bock	ζ
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${ m Analysis}$:	Mo, To	$_{ m tal}$	Analytical Method:			d: S (6010B		Prep Metho	od: S 3010A
QC Batch:	63221			Date	Analyzed:	20	09-09-03		Analyzed B	y: RR
Prep Batch:	53951			Samp	le Preparati	on: 20	09-09-03		Prepared B	y: KV
			SDL	$_{ m MQL}$	Method					
			Based	Based	Blank				MQL	MDL
Parameter		Flag	Result	Result	Result	Units	$\operatorname{Dilution}$	SDL	(Unadjusted)	(Unadjusted)
Total Molyb	denum		0.447	0.447	< 0.00119	$\mathrm{mg/L}$	1	0.00119	0.01	0.00119

Sample: 208954 - HLSF-0154-DRW-112-0809

Lacotratory.							
Analysis:	Na, Total		Ana	alytical Method:	S_{6010B}	Prep Method:	S 3010
QC Batch:	63355		Dat	e Analyzed:	2009-09-09	Analyzed By:	RR
Prep Batch:	53951		San	ple Preparation:	2009-09-03	Prepared By:	KV
		anı	MOT				

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Sodium		$\boldsymbol{5830}$	$\boldsymbol{5830}$	< 5.00	$\mathrm{mg/L}$	100	5.00	1	0.05

Sample: 208954 - HLSF-0154-DRW-112-0809

Report Date: September 18, 2009 Work Order: 9090211 Page Number: 12 of 80

HELSTF Diesel Spill Groundwater

Laboratory: Analysis: QC Batch: Prep Batch:	Lubbock Ni, Total 63221 53951		Γ	nalytical Met ate Analyzed ample Prepar	:	S 6010B 2009-09-03 2009-09-03		Analyzed E	Prep Method: S 3010A Analyzed By: RR Prepared By: KV		
		SDL	MQL	Method							
		Based	Based	Blank				MQL	MDL		
Parameter	Flag	Result	Result	Result	Units	$\operatorname{Dilution}$	SDL	(Unadjusted)	(Unadjusted)		
Total Nickel		0.780	0.780	< 0.00121	mg/L	1	0.00121	0.005	0.00121		

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Nitrate and Nitrite as N Analytical Method: SM 4500-NO3 EPrep Method: N/AQC Batch: 63485 Date Analyzed: 2009-09-11 Analyzed By: KVPrep Batch: 54188Sample Preparation: 2009-09-11 Prepared By: KV

SDLMQLMethod Based MQLMDLBased Blank Flag Parameter Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) 204 17.5 Nitrate and Nitrite as N 204< 17.5mg/L500 0.10.035

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: El Paso

O/G Analytical Method: Analysis: E 1664 Prep Method: N/AQC Batch: 63699 Date Analyzed: 2009-09-17 Analyzed By: JRPrep Batch: 54391 Sample Preparation: 2009-09-17 Prepared By: JR

SDLMQLMethod Based Based Blank MQLMDLFlag SDL Parameter Result Result Result Units Dilution (Unadjusted) (Unadjusted) 3.60 Oil and Grease < 3.60< 5.00< 3.60mg/L5 3.6

Sample: 208954 - HLSF-0154-DRW-112-0809

 $Laboratory : \quad Lubbock$

Analysis: P, Total Analytical Method: S 3010A S 6010B Prep Method: QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDLMQL MethodBased Based Blank MQLMDLParameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) Total Phosphorous 0.05500.0550< 0.00289 mg/L0.002890.0250.00289

Report Date: September 18, 2009 Work Order: 9090211 Page Number: 13 of 80

HELSTF Diesel Spill Groundwater

Laboratory: Lubbock

Analysis: Pb, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

 SDL MQL Method

MQLBased Based Blank MDL Dilution Parameter Result Result Result Units SDL(Unadjusted) Flag (Unadjusted) Total Lead < 0.00326 < 0.00500 < 0.00326 mg/L 0.00326 0.0050.00326

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: El Paso

Analysis: рΗ Analytical Method: SM 4500-H+Prep Method: N/AQC Batch: 63339 JGDate Analyzed: 2009-09-01 Analyzed By: Prep Batch: 54063 Sample Preparation: 2009-09-01 Prepared By: JR

RL

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Sb, Total Analysis: Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDL MQL Method Based Based Blank

Blank MQLMDL Flag SDLParameter Result Result Result Units Dilution (Unadjusted) (Unadjusted) 0.02 Total Antimony 0.0140< 0.0200< 0.00440mg/L0.004400.0044

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Se, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951KVSample Preparation: 2009-09-03 Prepared By:

SDL MQL Method

		Dased	Dased	ыанк				MCT	MDL
Parameter	Flag	Result	Result	Result	Units	$\operatorname{Dilution}$	SDL	(Unadjusted)	(Unadjusted)
Total Selenium		0.270	0.270	< 0.00508	$\mathrm{mg/L}$	1	0.00508	0.02	0.00508

MOT

MINT

HELSTF Diesel Spill Groundwater

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock Analysis: Semivolatiles Analytical Method: S 8270CPrep Method: S 3510C QC Batch: Analyzed By: 63261Date Analyzed: 2009-09-04 MN Prep Batch: 53996 Sample Preparation: 2009-09-02 Prepared By: MN

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Pyridine	J	0.00138	< 0.00490	< 0.000596	mg/L	0.98	0.000596	0.005	0.000608
N-Nitrosodimethylamine	U	< 0.000541	< 0.00490	< 0.000541	mg/L	0.98	0.000541	0.005	0.000552
2-Picoline	U	< 0.000400	< 0.00490	< 0.000400	mg/L	0.98	0.000400	0.005	0.000408
Methyl methanesulfonate	U	< 0.000343	< 0.00490	< 0.000343	mg/L	0.98	0.000343	0.005	0.00035
Ethyl methanesulfonate	U	< 0.000439	< 0.00490	< 0.000439	mg/L	0.98	0.000439	0.005	0.000448
Phenol	U	< 0.000499	< 0.00490	< 0.000499	mg/L	0.98	0.000499	0.005	0.000509
Aniline	U	< 0.000677	< 0.00490	< 0.000677	mg/L	0.98	0.000677	0.005	0.000691
bis(2-chloroethyl)ether	U	< 0.000431	< 0.00490	< 0.000431	mg/L	0.98	0.000431	0.005	0.00044
2-Chlorophenol	U	< 0.000526	< 0.00490	< 0.000526	mg/L	0.98	0.000526	0.005	0.000537
1,3-Dichlorobenzene (meta)	U	< 0.000432	< 0.00490	< 0.000432	mg/L	0.98	0.000432	0.005	0.000441
1,4-Dichlorobenzene (para)	U	< 0.000431	< 0.00490	< 0.000431	mg/L	0.98	0.000431	0.005	0.00044
Benzyl alcohol	U	< 0.000527	< 0.00490	< 0.000527	mg/L	0.98	0.000527	0.005	0.000538
1,2-Dichlorobenzene (ortho)	U	< 0.000434	< 0.00490	< 0.000434	mg/L	0.98	0.000434	0.005	0.000443
2-Methylphenol	U	< 0.000711	< 0.00490	< 0.000711	mg/L	0.98	0.000711	0.005	0.000726
bis(2-chloroisopropyl)ether	U	< 0.000493	< 0.00490	< 0.000493	mg/L	0.98	0.000493	0.005	0.000503
4-Methylphenol / 3-Methylphenol		< 0.000502	< 0.00490	< 0.000502	mg/L	0.98	0.000502	0.005	0.000512
N-Nitrosodi-n-propylamine	U	< 0.000717	< 0.00490	< 0.000717	mg/L	0.98	0.000717	0.005	0.000732
${\it Hexachloroethane}$	U	< 0.000497	< 0.00490	< 0.000497	mg/L	0.98	0.000497	0.005	0.000507
$egin{array}{c} Acetophenone \end{array}$	U	< 0.000416	< 0.00490	< 0.000416	mg/L	0.98	0.000416	0.005	0.000424
Nitrobenzene	U	< 0.000456	< 0.00490	< 0.000456	mg/L	0.98	0.000456	0.005	0.000465
N-Nitrosopiperidine	U	< 0.000434	< 0.00490	< 0.000434	mg/L	0.98	0.000434	0.005	0.000443
Isophorone	U	< 0.000607	< 0.00490	< 0.000607	mg/L	0.98	0.000607	0.005	0.000619
2-Nitrophenol	U	< 0.000398	< 0.00490	< 0.000398	mg/L	0.98	0.000398	0.005	0.000406
2,4-Dimethylphenol	U	< 0.000467	< 0.00490	< 0.000467	mg/L	0.98	0.000467	0.005	0.000477
bis(2-chloroethoxy) methane	U	< 0.000423	< 0.00490	< 0.000423	mg/L	0.98	0.000423	0.005	0.000432
2,4-Dichlorophenol	U	< 0.000392	< 0.00490	< 0.000392	mg/L	0.98	0.000392	0.005	0.0004
1,2,4-Trichlorobenzene	U	< 0.000396	< 0.00490	< 0.000396	mg/L	0.98	0.000396	0.005	0.000404
Benzoic acid	U	< 0.00160	< 0.00490	< 0.00160	mg/L	0.98	0.00160	0.005	0.00163
Naphthalene	U	< 0.000479	< 0.00490	< 0.000479	mg/L	0.98	0.000479	0.005	0.000489
a,a-Dimethylphenethylamine	U	< 0.00126	< 0.00490	< 0.00126	mg/L	0.98	0.00126	0.005	0.00129
4-Chloroaniline	U	< 0.000370	< 0.00490	< 0.000370	mg/L	0.98	0.000370	0.005	0.000378
2,6-Dichlorophenol	U	< 0.000474	< 0.00980	< 0.000474	mg/L	0.98	0.000474	0.01	0.000484
Hexachlorobutadiene	U	< 0.000507	< 0.00490	< 0.000507	mg/L	0.98	0.000507	0.005	0.000517
N-Nitroso-di-n-butylamine	U	< 0.000643	< 0.00490	< 0.000643	mg/L	0.98	0.000643	0.005	0.000656
4-Chloro-3-methylphenol	U	< 0.000512	< 0.00490	< 0.000512	mg/L	0.98	0.000512	0.005	0.000522
2-Methylnaphthalene	U	< 0.000414	< 0.00490	< 0.000414	mg/L	0.98	0.000414	0.005	0.000423
$1\hbox{-}Methylnaphthalene$	U	< 0.000485	< 0.00490	< 0.000485	mg/L	0.98	0.000485	0.005	0.000495
1,2,4,5-Tetrachlorobenzene	U	< 0.000600	< 0.00490	< 0.000600	mg/L	0.98	0.000600	0.005	0.000612
${\it Hexachlorocyclopentadiene}$	U	< 0.000547	< 0.00490	< 0.000547	mg/L	0.98	0.000547	0.005	0.000558
2,4,6-Trichlorophenol	U	< 0.000778	< 0.00980	< 0.000778	mg/L	0.98	0.000778	0.01	0.000794

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sample 208954 continued . . .

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
2,4,5-Trichlorophenol	U	< 0.000817	< 0.00490	< 0.000817	mg/L	0.98	0.000817	0.005	0.000834
$2 ext{-Chloronaphthalene}$	U	< 0.000408	< 0.00490	< 0.000408	mg/L	0.98	0.000408	0.005	0.000416
$1 ext{-Chloronaphthalene}$	U	< 0.000466	< 0.00490	< 0.000466	mg/L	0.98	0.000466	0.005	0.000476
2-Nitroaniline	U	< 0.000745	< 0.00490	< 0.000745	mg/L	0.98	0.000745	0.005	0.00076
${ m Dimethyl}{ m phthalate}$	U	< 0.000630	< 0.00490	< 0.000630	mg/L	0.98	0.000630	0.005	0.000643
${ m Acenaphthylene}$	U	< 0.000574	< 0.00490	< 0.000574	mg/L	0.98	0.000574	0.005	0.000586
2,6-Dinitrotoluene	U	< 0.000627	< 0.00490	< 0.000627	mg/L	0.98	0.000627	0.005	0.00064
3-Nitroaniline	U	< 0.000706	< 0.00490	< 0.000706	mg/L	0.98	0.000706	0.005	0.000721
${ m Acenaphthene}$	U	< 0.000414	< 0.00490	< 0.000414	mg/L	0.98	0.000414	0.005	0.000423
2,4-Dinitrophenol	U	< 0.000216	< 0.00490	< 0.000216	mg/L	0.98	0.000216	0.005	0.00022
Dibenzofuran	U	< 0.000400	< 0.00490	< 0.000400	mg/L	0.98	0.000400	0.005	0.000408
${ m Pentachlorobenzene}$	U	< 0.000560	< 0.00490	< 0.000560	mg/L	0.98	0.000560	0.005	0.000571
4-Nitrophenol	U	< 0.00181	< 0.0245	< 0.00181	mg/L	0.98	0.00181	0.025	0.00185
2,4-Dinitrotoluene	U	< 0.000893	< 0.00490	< 0.000893	mg/L	0.98	0.000893	0.005	0.000911
1-Naphthylamine	U	< 0.000674	< 0.00490	< 0.000674	mg/L	0.98	0.000674	0.005	0.000688
2,3,4,6-Tetrachlorophenol	U	< 0.000554	< 0.00980	< 0.000554	mg/L	0.98	0.000554	0.01	0.000565
2-Naphthylamine	U	< 0.000685	< 0.00490	< 0.000685	mg/L	0.98	0.000685	0.005	0.000699
Fluorene	U	< 0.000635	< 0.00490	< 0.000635	mg/L	0.98	0.000635	0.005	0.000648
4-Chlorophenyl-phenylether	U	< 0.000607	< 0.00490	< 0.000607	mg/L	0.98	0.000607	0.005	0.000619
${ m Diet}$ hylphthalate	U	< 0.000811	< 0.00490	< 0.000811	mg/L	0.98	0.000811	0.005	0.000828
4-Nitroaniline	U	< 0.000688	< 0.00490	< 0.000688	mg/L	0.98	0.000688	0.005	0.000702
Diphenylhydrazine	U	< 0.000644	< 0.00490	< 0.000644	mg/L	0.98	0.000644	0.005	0.000657
4,6-Dinitro- 2 -methylphenol	U	< 0.00194	< 0.00490	< 0.00194	mg/L	0.98	0.00194	0.005	0.00198
Diphenylamine	U	< 0.000431	< 0.00490	< 0.000431	mg/L	0.98	0.000431	0.005	0.00044
4-Bromophenyl-phenylether	U	< 0.000539	< 0.00490	< 0.000539	mg/L	0.98	0.000539	0.005	0.00055
Phenacetin	U	< 0.000593	< 0.00490	< 0.000593	mg/L	0.98	0.000593	0.005	0.000605
${ m Hexachlorobenzene}$	U	< 0.000496	< 0.00490	< 0.000496	mg/L	0.98	0.000496	0.005	0.000506
4-Aminobiphenyl	U	< 0.000516	< 0.00490	< 0.000516	mg/L	0.98	0.000516	0.005	0.000527
${ m Pentachlorophenol}$	U	< 0.000426	< 0.00980	< 0.000426	mg/L	0.98	0.000426	0.01	0.000435
${ m Anthracene}$	U	< 0.000419	< 0.00490	< 0.000419	mg/L	0.98	0.000419	0.005	0.000428
${\bf Pentachloronitrobenzene}$	U	< 0.000400	< 0.00490	< 0.000400	mg/L	0.98	0.000400	0.005	0.000408
Pronamide	U	< 0.000466	< 0.00490	< 0.000466	mg/L	0.98	0.000466	0.005	0.000476
${ m Phenanthrene}$	U	< 0.000537	< 0.00490	< 0.000537	mg/L	0.98	0.000537	0.005	0.000548
${ m Di} ext{-n-butylpht}$ halate	U	< 0.000473	< 0.00490	< 0.000473	mg/L	0.98	0.000473	0.005	0.000483
Fluoranthene	U	< 0.000619	< 0.00490	< 0.000619	mg/L	0.98	0.000619	0.005	0.000632
Benzidine	U	< 0.00233	< 0.0245	< 0.00233	mg/L	0.98	0.00233	0.025	0.00238
Pyrene	U	< 0.000708	< 0.00490	< 0.000708	mg/L	0.98	0.000708	0.005	0.000723
${\rm p\text{-}Dimethylaminoazobenzene}$	U	< 0.000884	< 0.00490	< 0.000884	mg/L	0.98	0.000884	0.005	0.000902
${f Butylbenzylphthalate}$	U	< 0.000436	< 0.00490	< 0.000436	mg/L	0.98	0.000436	0.005	0.000445
$\operatorname{Benzo}(\operatorname{a}) \operatorname{anthracene}$	U	< 0.000516	< 0.00490	< 0.000516	mg/L	0.98	0.000516	0.005	0.000527
3,3-Dichlorobenzidine	U	< 0.00116	< 0.00490	< 0.00116	mg/L	0.98	0.00116	0.005	0.00118
Chrysene	U	< 0.000625	< 0.00490	< 0.000625	mg/L	0.98	0.000625	0.005	0.000638
bis(2-ethylhexyl)phthalate	U	< 0.000550	< 0.00490	< 0.000550	mg/L	0.98	0.000550	0.005	0.000561
Di-n-octylphthalate	U			< 0.00114		0.98	0.00114	0.005	0.00116
Benzo(b)fluoranthene	U	< 0.000861	< 0.00490	< 0.000861	mg/L	0.98	0.000861	0.005	0.000879

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${\it Work~Order:~9090211} \\ {\it HELSTF~Diesel~Spill~Groundwater}$

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sample 208954 continued . . .

		SDL	MQL	$\operatorname{Met} \operatorname{hod}$				
		Based	Based	Blank			MQL	MDL
Parameter	Flag	Result	Result	Result Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
$\overline{\mathrm{Benzo}(\mathtt{k})}$ fluoranthene	U	< 0.000828	< 0.00490	$< 0.000828 \mathrm{mg/L}$	0.98	0.000828	0.005	0.000845
7,12-Dimethylbenz(a)anthracene	U	< 0.00100	< 0.00490	$< 0.00100 \mathrm{mg/L}$	0.98	0.00100	0.005	0.00102
$\mathrm{Benzo}(\mathrm{a})\mathrm{pyrene}$	U	< 0.00164	< 0.00490	< 0.00164 mg/L	0.98	0.00164	0.005	0.00167
3-Methylcholanthrene	U	< 0.000890	< 0.00490	$< 0.000890 \; \mathrm{mg/L}$	0.98	0.000890	0.005	0.000908
${ m Dibenzo(a,j)}$ acridine	U	< 0.00126	< 0.00490	< 0.00126 mg/L	0.98	0.00126	0.005	0.00129
Indeno(1,2,3-cd)pyrene	U	< 0.000845	< 0.00490	< 0.000845 mg/L	0.98	0.000845	0.005	0.000862
${ m Dibenzo(a,h)}$ anthracene	U	< 0.000793	< 0.00490	< 0.000793 mg/L	0.98	0.000793	0.005	0.000809
$\mathrm{Benzo}(\mathrm{g,h,i})$ perylene	U	< 0.000930	< 0.00490	${<}0.000930\:\mathrm{mg/L}$	0.98	0.000930	0.005	0.000949

					Spike	$\operatorname{Percent}$	Recovery
Surrogate	Flag	Result	Units	Dilution	${f Amount}$	Recovery	Limits
2-Fluorophenol		0.0276	$_{ m mg/L}$	0.98	0.0800	34	10 - 53.1
${ m Phenol-d5}$		0.0191	${ m mg/L}$	0.98	0.0800	24	10 - 36.9
${ m Nitrobenzene-d5}$		0.0447	${ m mg/L}$	0.98	0.0800	56	23.8 - 108
2-Fluorobiphenyl		0.0439	${ m mg/L}$	0.98	0.0800	55	15.9 - 127
2,4,6-Tribromophenol		0.0594	${ m mg/L}$	0.98	0.0800	74	10 - 123
${ m Terphenyl-d14}$		0.0520	$\mathrm{mg/L}$	0.98	0.0800	65	17.2 - 160

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: El Paso

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Sulfate		11600	11600	< 504	mg/L	1000	504	1.33	0.5038

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: El Paso

Analysis: TDS Analytical Method: SM 2540C Prep Method: N/A QC Batch: 63473 Date Analyzed: 2009-09-03 Analyzed By: MD Prep Batch: 54173 Sample Preparation: 2009-09-03 Prepared By: MD

		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Dissolved Solids		29200	29200	< 5.00	mg/L	1	5.00		5

HELSTF Diesel Spill Groundwater

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: TKN Analytical Method: E 351.3 Prep Method: N/AQC Batch: 63441 Date Analyzed: 2009-09-09 Analyzed By: AHPrep Batch: 54150 Sample Preparation: 2009-09-09 Prepared By: AH

SDL MQL Method

MQLBased Based Blank MDL Parameter Result Result Result Units Dilution SDL(Unadjusted) Flag (Unadjusted) Total Kjeldahl Nitrogen - N < 2.45 <10.0 < 2.45mg/L 2.4510 2.45 1

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Tl, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 2009-09-03 Analyzed By: 63221 Date Analyzed: RRPrep Batch: 53951 Sample Preparation: 2009-09-03 Prepared By: KV

SDL MQL Method

MQLMDL Based Based Blank SDLDilution Parameter Flag Result Result Result Units (Unadjusted) (Unadjusted) Total Thallium < 0.00488 < 0.0500 < 0.00488mg/L1 0.00488 0.050.00488

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: TOC Analytical Method: SM 5310CPrep Method: N/AQC Batch: 63340 Date Analyzed: 2009-09-08 Analyzed By: KVPrep Batch: 54065Sample Preparation: 2009-09-08 Prepared By: KV

 $\operatorname{SDL} \quad \operatorname{MQL} \quad \operatorname{Method}$

Based MQLMDL Based Blank Parameter Result Result Result Dilution SDLFlag Units (Unadjusted) (Unadjusted) Total Organic Carbon 1.181.18< 0.401mg/L0.4010.401

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Total Cyanide Analytical Method: SM 4500-CN C,E Prep Method: N/AQC Batch: AH63391 Date Analyzed: 2009-09-07 Analyzed By: Prep Batch: 54106Sample Preparation: 2009-09-07 Prepared By: AH

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sample 208954 continued . . .

Parameter	Flag	$\begin{array}{c} \mathrm{SDL} \\ \mathrm{Based} \\ \mathrm{Result} \end{array}$	$egin{array}{l} \mathrm{MQL} \\ \mathrm{Based} \\ \mathrm{Result} \end{array}$	Method Blank Result	${ m Units}$	Dilution	SDL	$egin{aligned} \mathbf{MQL} \ & (\mathtt{Unadjusted}) \end{aligned}$	MDL (Unadjusted)
		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Cyanide	U	< 0.0110	< 0.0150	< 0.0110	mg/L	1	0.0110	0.015	0.011

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: TPH DRO Analytical Method: Prep Method: N/A Mod. 8015B

QC Batch: 63307 Date Analyzed: 2009-09-07Analyzed By: Prep Batch: Sample Preparation: 540352009-09-04 Prepared By:

SDLMQLMethod Based Based Blank MQLMDLSDLParameter Flag Result Result Result Units Dilution (Unadjusted) (Unadjusted) DRO < 0.8760.876 < 0.876< 5.00mg/L5 0.876

					Spike	$\operatorname{Percent}$	$\operatorname{Recovery}$
Surrogate	Flag	Result	Units	Dilution	${f Amount}$	Recovery	Limits
n-Triacontane		10.4	m mg/L	1	10.0	104	57.3 - 151

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: TPH GRO Analytical Method: S 8015BPrep Method: S 5030B QC Batch: 63239 Date Analyzed: 2009-09-03 Analyzed By: ERPrep Batch: Sample Preparation: 2009-09-03 53976 Prepared By: ER

SDLMQLMethod Based Based Blank MQLMDLParameter Flag Result Result Result Units Dilution SDL(Unadjusted) (Unadjusted) GRO < 0.152< 0.200 < 0.152mg/L0.1520.2 0.1521

Surrogate	Flag	Result	Units	Dilution	$\begin{array}{c} {\rm Spike} \\ {\rm Amount} \end{array}$	Percent Recovery	$egin{array}{l} { m Recovery} \\ { m Limits} \end{array}$
Trifluorotoluene (TFT)		0.105	mg/L	1	0.100	105	70.8 - 112
4-Bromofluorobenzene (4-BFB)		0.105	${ m mg/L}$	1	0.100	105	80 - 109

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Report Date: September 18, 2009 Work Order: 9090211 Page Number: 19 of 80

HELSTF Diesel Spill Groundwater

Analysis: QC Batch: Prep Batch:	V, Total 63221 53951		Date	lytical Metho e Analyzed: ple Preparati	200	010B 9-09-03 9-09-03		Prep Metho Analyzed B Prepared B	By: RR
		SDL	MQL	Method					
		Based	Based	Blank				MQL	MDL
Parameter	Flag	Result	Result	Result	Units	Dilution	SDL	(Unadjusted)	(Unadjusted)
Total Vanadi	um	0.0100	0.0100	< 0.000426	mg/L	1	0.000426	0.005	0.000426

Sample: 208954 - HLSF-0154-DRW-112-0809

Laboratory: Lubbock

Analysis: Zn, Total Analytical Method: S 6010B Prep Method: S 3010A QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RRPrep Batch: 53951Sample Preparation: 2009-09-03 Prepared By: KV

SDL MQLMethod MQLMDL Based Based Blank Result SDLParameter Flag Result Result Units Dilution (Unadjusted) (Unadjusted) Total Zinc < 0.000465 < 0.00500 < 0.000465 mg/L0.0004650.0050.000465

Method Blank (1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

Method Blank (1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

Method Blank (1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

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		-		
Parameter	Flag	Result	Units	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$
Total Arsenic	1 148	<0.00448	mg/L	0.00448
Method Blank (1)				
QC Batch: 63221		Date Analyzed: 2009-09-03		Analyzed By: RR
Prep Batch: 53951		QC Preparation: 2009-09-03		Prepared By: KV
				Reporting
Parameter	Flag	Result	Units	Limits
Total Barium		< 0.00105	m mg/L	0.00105
Method Blank (1)				
QC Batch: 63221		Date Analyzed: 2009-09-03		Analyzed By: RR
Prep Batch: 53951		QC Preparation: 2009-09-03		Prepared By: KV
				Reporting
Parameter	Flag	Result	Units	Limits
Total Beryllium		< 0.000450	mg/L	0.00045
Method Blank (1)				
QC Batch: 63221		Date Analyzed: 2009-09-03		Analyzed By: RR
Prep Batch: 53951		QC Preparation: 2009-09-03		Prepared By: KV
				Reporting
Parameter	Flag	Result	Units	Limits
Total Cadmium		< 0.000303	m mg/L	0.000303
Method Blank (1)				
QC Batch: 63221		Date Analyzed: 2009-09-03		Analyzed By: RR
Prep Batch: 53951		QC Preparation: 2009-09-03		Prepared By: KV
		_		Reporting
Parameter	Flag	Result	Units	Limits
Total Cobalt		< 0.000822	${ m mg/L}$	0.000822

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HELSTF Diesel Spill Groundwater

		HELSTF Diesel Spill Groundw	ater	
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	Result	Units	Reporting Limits
Total Chromium		< 0.000583	m mg/L	0.000583
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	Result	Units	Reporting Limits
Total Copper		< 0.000843	m mg/L	0.000843
Method Blank (1) QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	m Result	Units	Reporting Limits
Total Iron		< 0.000872	m mg/L	0.000872
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter Total Manganese	Flag	Result <0.000305	$\begin{array}{c} \rm Units \\ \rm mg/L \end{array}$	Reporting Limits 0.000305
Total Manganese		<0.000909	mg/ L	0.000303
Method Blank (1)				

Date Analyzed:

QC Preparation: 2009-09-03

2009-09-03

Analyzed By: RR

Prepared By: KV

QC Batch:

Prep Batch: 53951

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Parameter	Flag	Result	Units	Reporting Limits
Total Molybdenum	9	< 0.00119	m mg/L	0.00119
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter Total Nickel	Flag	Result <0.00121	$\frac{\rm Units}{\rm mg/L}$	Reporting Limits 0.00121
Method Blank (1) QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter Total Phosphorous	Flag	Result <0.00289	$\begin{array}{c} \rm Units \\ \rm mg/L \end{array}$	Reporting Limits 0.00289
Method Blank (1) QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter Total Lead	Flag	Result <0.00326	$\frac{\rm Units}{\rm mg/L}$	Reporting Limits 0.00326
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter Total Antimony	Flag	Result < 0.00440	$rac{ m Units}{ m mg/L}$	Reporting Limits 0.0044

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HELSTF Diesel Spill Groundwater

-		HELSTF Diesel Spill Groundw	auci	
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	Result	Units	Reporting Limits
Total Selenium		< 0.00508	m mg/L	0.00508
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	Result	Units	$rac{ ext{Reporting}}{ ext{Limits}}$
Total Thallium		< 0.00488	m mg/L	0.00488
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	Result	${ m Units}$	$rac{ ext{Reporting}}{ ext{Limits}}$
Total Vanadium		< 0.000426	m mg/L	0.000426
Method Blank (1)				
QC Batch: 63221 Prep Batch: 53951		Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
Parameter	Flag	Result	${ m Units}$	Reporting Limits
Total Zinc		< 0.000465	$\mathrm{mg/L}$	0.000465

Date Analyzed:

QC Preparation: 2009-09-03

2009 - 09 - 03

Analyzed By: TP

Prepared By: TP

QC Batch:

Prep Batch: 53960

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				Reporting
Parameter	Flag	Result	Units	Limits
Total Mercury		< 0.0000329	$_{ m mg/L}$	3.29e-05

Method Blank (1)

QC Batch: Date Analyzed: Analyzed By: ER 63239 2009-09-03 Prep Batch: 53976 QC Preparation: 2009-09-03 Prepared By: ER

							Reporting
Parameter	Flag		Result		Units		Limits
GRO			< 0.152		${ m mg/L}$		0.152
					Spike	Percent	Recovery
Surrogate	Flag	Result	Units	$\operatorname{Dilution}$	${ m Amount}$	Recovery	Limits
Trifluorotoluene (TFT)		0.103	m mg/L	1	0.100	103	70.8 - 112
4-Bromofluorobenzene (4-BFB	3)	0.100	$_{ m mg/L}$	1	0.100	100	80 - 109

Method Blank (1)

QC Batch: 63261Date Analyzed: 2009-09-04 Analyzed By: MN Prep Batch: 53996 QC Preparation: 2009-09-02 Prepared By: MN

				Reporting
Parameter	Flag	Result	Units	Limits
Pyridine		< 0.000608	m mg/L	0.000608
N-Nitrosodimethylamine		< 0.000552	m mg/L	0.000552
2-Picoline		< 0.000408	m mg/L	0.000408
Methyl methanesulfonate		< 0.000350	m mg/L	0.00035
Ethyl methanesulfonate		< 0.000448	m mg/L	0.000448
Phenol		< 0.000509	m mg/L	0.000509
Aniline		< 0.000691	m mg/L	0.000691
bis(2-chloroethyl)ether		< 0.000440	m mg/L	0.00044
2-Chlorophenol		< 0.000537	m mg/L	0.000537
1,3-Dichlorobenzene (meta)		< 0.000441	m mg/L	0.000441
1,4-Dichlorobenzene (para)		< 0.000440	m mg/L	0.00044
Benzyl alcohol		< 0.000538	m mg/L	0.000538
1,2-Dichlorobenzene (ortho)		< 0.000443	m mg/L	0.000443
2-Methylphenol		< 0.000726	m mg/L	0.000726
bis(2-chloroisopropyl) ether		< 0.000503	m mg/L	0.000503
4-Methylphenol / 3-Methylphenol		< 0.000512	m mg/L	0.000512
N-Nitrosodi-n-propylamine		< 0.000732	m mg/L	0.000732
${\bf Hexachloroethane}$		< 0.000507	m mg/L	0.000507
$egin{array}{c} Acetophenone \end{array}$		< 0.000424	m mg/L	0.000424
Nitrobenzene		< 0.000465	m mg/L	0.000465
N-Nitrosopiperidine		< 0.000443	m mg/L	0.000443
Isophorone		< 0.000619	m mg/L	0.000619

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Parameter	Flag	Result	Units	Limits
2-Nitrophenol		< 0.000406	m mg/L	0.000406
2,4-Dimethylphenol		< 0.000477	m mg/L	0.000477
bis(2-chloroethoxy) methane		< 0.000432	m mg/L	0.000432
2,4-Dichlorophenol		< 0.000400	m mg/L	0.0004
1,2,4-Trichlorobenzene		< 0.000404	m mg/L	0.000404
Benzoic acid		< 0.00163	m mg/L	0.00163
Naphthalene		< 0.000489	m mg/L	0.000489
${\rm a,a-Dimethyl phenethyl amine}$		< 0.00129	m mg/L	0.00129
4-Chloroaniline		< 0.000378	m mg/L	0.000378
2,6-Dichlorophenol		< 0.000484	m mg/L	0.000484
Hexachlorobutadiene		< 0.000517	m mg/L	0.000517
N-Nitroso-di-n-butylamine		< 0.000656	m mg/L	0.000656
4-Chloro-3-methylphenol		< 0.000522	m mg/L	0.000522
2-Methylnaphthalene		< 0.000423	m mg/L	0.000423
$1 ext{-Methylnaphthalene}$		< 0.000495	m mg/L	0.000495
1,2,4,5-Tetrachlorobenzene		< 0.000612	m mg/L	0.000612
${\it Hexachlorocyclopenta}$ diene		< 0.000558	m mg/L	0.000558
2,4,6-Trichlorophenol		< 0.000794	m mg/L	0.000794
2,4,5-Trichlorophenol		< 0.000834	m mg/L	0.000834
$2 ext{-Chloronaphthalene}$		< 0.000416	m mg/L	0.000416
$1 ext{-}\mathrm{Chloronaphthalene}$		< 0.000476	m mg/L	0.000476
2-Nitroaniline		< 0.000760	m mg/L	0.00076
${\bf Dimethylphthalate}$		< 0.000643	m mg/L	0.000643
${f Acenaphthylene}$		< 0.000586	m mg/L	0.000586
2,6-Dinitrotoluene		< 0.000640	m mg/L	0.00064
3-Nitroaniline		< 0.000721	m mg/L	0.000721
${ m Acenapht}$ hene		< 0.000423	m mg/L	0.000423
2,4-Dinitrophenol		< 0.000220	m mg/L	0.00022
Dibenzofuran		< 0.000408	m mg/L	0.000408
Pentachlorobenzene		< 0.000571	m mg/L	0.000571
4-Nitrophenol		< 0.00185	m mg/L	0.00185
2,4-Dinitrotoluene		< 0.000911	m mg/L	0.000911
1-Naphthylamine		< 0.000688	m mg/L	0.000688
2,3,4,6-Tetrachlorophenol		< 0.000565	m mg/L	0.000565
2-Naphthylamine		< 0.000699	m mg/L	0.000699
Fluorene		< 0.000648	m mg/L	0.000648
4-Chlorophenyl-phenylether		< 0.000619	m mg/L	0.000619
Diethylphthalate		< 0.000828	m mg/L	0.000828
4-Nitroaniline		< 0.000702	m mg/L	0.000702
Diphenylhydrazine		< 0.000657	m mg/L	0.000657
4,6-Dinitro- 2 -methylphenol		< 0.00198	m mg/L	0.00198
Diphenylamine		< 0.000440	m mg/L	0.00044
4-Bromophenyl-phenylether		< 0.000550	m mg/L	0.00055
Phenacetin		< 0.000605	m mg/L	0.000605
Hexachlorobenzene		< 0.000506	m mg/L	0.000506
4-Aminobiphenyl		< 0.000527	m mg/L	0.000527
Pentachlorophenol		< 0.000435	m mg/L	0.000435

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method	blank	continued		

				Reporting
Parameter	Flag	Result	Units	Limits
Anthracene		< 0.000428	m mg/L	0.000428
${\bf Pentach loronitrobenzene}$		< 0.000408	m mg/L	0.000408
Pronamide		< 0.000476	m mg/L	0.000476
Phenanthrene		< 0.000548	m mg/L	0.000548
Di-n-butylphthalate		< 0.000483	m mg/L	0.000483
Fluoranthene		< 0.000632	m mg/L	0.000632
Benzidine		< 0.00238	m mg/L	0.00238
Pyrene		< 0.000723	m mg/L	0.000723
${\bf p-Dimethylaminoazobenzene}$		< 0.000902	m mg/L	0.000902
${\bf Butylbenzylphthalate}$		< 0.000445	m mg/L	0.000445
$\operatorname{Benzo}(\operatorname{a}) \operatorname{anthracene}$		< 0.000527	m mg/L	0.000527
3,3-Dichlorobenzidine		< 0.00118	m mg/L	0.00118
Chrysene		< 0.000638	m mg/L	0.000638
${ m bis}(2{ m -ethylhexyl}){ m phthalate}$		< 0.000561	m mg/L	0.000561
Di-n-octylphthalate		< 0.00116	m mg/L	0.00116
$\operatorname{Benzo}(b)$ fluoranthene		< 0.000879	m mg/L	0.000879
$\operatorname{Benzo}(k)$ fluoranthene		< 0.000845	m mg/L	0.000845
7,12-Dimethylbenz(a) anthracene		< 0.00102	m mg/L	0.00102
$\mathrm{Benzo}(\mathrm{a})\mathrm{pyrene}$		< 0.00167	m mg/L	0.00167
3-Methylcholanthrene		< 0.000908	m mg/L	0.000908
${ m Dibenzo(a,j)}$ acridine		< 0.00129	m mg/L	0.00129
Indeno(1,2,3-cd)pyrene		< 0.000862	m mg/L	0.000862
${ m Dibenzo(a,h)}$ anthracene		< 0.000809	m mg/L	0.000809
Benzo(g,h,i)perylene		< 0.000949	m mg/L	0.000949

					Spike	Percent	Recovery
Surrogate	Flag	Result	Units	$\operatorname{Dilution}$	${f Amount}$	Recovery	Limits
2-Fluorophenol		0.0250	${ m mg/L}$	1	0.0800	31	10 - 53.1
${ m Phenol-d5}$		0.0158	${ m mg/L}$	1	0.0800	20	10 - 36.9
${ m Nitrobenzene-d5}$		0.0378	${ m mg/L}$	1	0.0800	47	23.8 - 108
2-Fluorobiphenyl		0.0396	${ m mg/L}$	1	0.0800	50	15.9 - 127
2,4,6-Tribromophenol		0.0493	$_{ m mg/L}$	1	0.0800	62	10 - 123
Terphenyl-d14		0.0437	$_{ m mg/L}$	1	0.0800	55	17.2 - 160

Method Blank (1)

QC Batch: 63307 Date Analyzed: 2009-09-07 Analyzed By: Prep Batch: 54035 QC Preparation: 2009-09-04 Prepared By:

				Reporting
Parameter	Flag	Result	${ m Units}$	Limits
DRO		< 0.876	m mg/L	0.876

					Spike	$\operatorname{Percent}$	Recovery
$\mathbf{Surrogate}$	Flag	Result	Units	Dilution	${f Amount}$	Recovery	Limits
n-Triacontane		8.25	$\mathrm{mg/L}$	1	10.0	82	57.3 - 151

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HELSTF Diesel Spill Groundwater

	1122	orr Broson Spin Ground.		
Method Blank (1)				
QC Batch: 63327 Prep Batch: 54055	Date Ar	nalyzed: 2009-09-01 paration: 2009-09-01		Analyzed By: MD Prepared By: MD
Frep Datch: 54055	QC Fre	paration: 2009-09-01		Frepared by: MD
Parameter	Flor	Result	Units	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$
Hexavalent Chromium	Flag	<0.00594	mg/L	0.00594
			- Oi	
Method Blank (1)				
QC Batch: 63340	Date Ar			Analyzed By: KV
Prep Batch: 54065	QC Pre	paration: 2009-09-08		Prepared By: KV
				Reporting
Parameter	Flag	Result	Units	Limits
Total Organic Carbon		0.848	mg/L	0.401
Method Blank (1)				
` ,	T	1 1 2000 00 00		4 1 1D DD
QC Batch: 63355 Prep Batch: 53951	Date Ar QC Pre	nalyzed: 2009-09-09 paration: 2009-09-03		Analyzed By: RR Prepared By: KV
Trop Baccin 00001	Q 2 2 2	paraoisii. 2000 00 00		-
Parameter	Flag	Result	Units	Reporting Limits
Total Calcium	Flag	<0.117	mg/L	0.117
			01	
Method Blank (1)				
QC Batch: 63355	Date Ar	· ·		Analyzed By: RR
Prep Batch: 53951	QC Pre	paration: 2009-09-03		Prepared By: KV
				Reporting
Parameter	Flag	Result	Units	Limits
Total Potassium		< 0.172	m mg/L	0.172
Method Blank (1)				
QC Batch: 63355	Date Aı	nalyzed: 2009-09-09		Analyzed By: RR
Prep Batch: 53951		paration: 2009-09-03		Prepared By: KV
				Reporting
Parameter	Flag	Result	Units	Limits
Total Magnesium		< 0.160	m mg/L	0.16

		HELSTF Diesel Spill Ground	water	
Method Blank (1)				
QC Batch: 63355 Prep Batch: 53951		e Analyzed: 2009-09-09 Preparation: 2009-09-03		Analyzed By: RR Prepared By: KV
				Reporting
Parameter	Flag	Result	Units	Limits
Total Sodium		< 0.0500	m mg/L	0.05
Method Blank (1)				
QC Batch: 63370	Dat	se Analyzed: 2009-09-05		Analyzed By: AH
Prep Batch: 54092		Preparation: 2009-09-05		Prepared By: AH
				T
Parameter	Flag	Result	Units	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$
Ammonia-N	Tag	<0.353	mg/L	0.353
		(0.000	8/ 12	0.000
Method Blank (1)				
QC Batch: 63391	Dat	se Analyzed: 2009-09-07		Analyzed By: AH
Prep Batch: 54106		Preparation: 2009-09-07		Prepared By: AH
D	-	D	T T 1.	Reporting
Parameter Total Cyanide	Flag	Result < 0.0110	$\frac{\rm Units}{\rm mg/L}$	Limits 0.011
		X0.0220	6/	0.022
Method Blank (1)				
QC Batch: 63425 Prep Batch: 54137		te Analyzed: 2009-09-10 Preparation: 2009-09-04		Analyzed By: DS Prepared By: DS
				Reporting
Parameter	Flag	Result	Units	Limits
HMX		< 0.123	$\mu \mathrm{g/L}$	0.123
RDX		< 0.298	$\mu \mathrm{g}/\mathrm{L}$	0.298
1,3,5-Trinitrobenzene 1,3-Dinitrobenzene		< 0.339 < 0.389	$\mu \mathrm{g}/\mathrm{L}$	$0.339 \\ 0.389$
Nitrobenzene		<0.389 <0.379	$\mu \mathrm{g}/\mathrm{L}$	0.389 0.379
Tetryl		< 0.413	$\mu \mathrm{g}/\mathrm{L}$	0.413
TNT		<0.464	$ m \mu g/L \ \mu g/L$	0.413 0.464
4-Amino-DNT		< 0.319	$ m \mu g/L \ \mu g/L$	0.404 0.319
2-Amino-DNT		< 0.391	$_{ m \mu g/L}^{ m \mu g/L}$	0.319 0.391
2,6-DNT		< 0.323	$_{ m \mu g/L}^{ m \mu g/L}$	0.323
2,4-DNT		< 0.366	$_{ m \mu g/L}^{ m \mu g/L}$	0.366
2-NT		< 0.379	$\mu_{ m g/L}$	0.379
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Report Date: September 18, 2		Work Order: 90902 STF Diesel Spill Grou	Page Number: 29 of 80		
method blank continued					
Parameter	Flag	Result	Units	$egin{array}{c} ext{Reporting} \ ext{Limits} \end{array}$	
4-NT	1 148	< 0.398	$\frac{\mu g/L}$	0.398	
3-NT		< 0.346	$\mu \mathrm{g}/\mathrm{L}$	0.346	
			Spike	Percent Recovery	
Surrogate Flag		nits Dilution		Recovery Limits	
1,2-Dinitrobenzene	2.40 μ	g/L 1	2.50	96 19.8 - 160	
Method Blank (1)					
QC Batch: 63441 Prep Batch: 54150		nalyzed: 2009-09-09 eparation: 2009-09-09		Analyzed By: AH Prepared By: AH	
D	T)	D 1	TT '	Reporting	
Parameter Total Kjeldahl Nitrogen - N	Flag	Result <2.45	$\frac{\text{Units}}{\text{mg/L}}$	Limits 2.45	
QC Batch: 63461 Prep Batch: 54153		nalyzed: 2009-09-11 eparation: 2009-09-11 Result	Units	Analyzed By: RR Prepared By: KV Reporting Limits	
Dissolved Chromium		< 0.000583	m mg/L	0.000583	
Method Blank (1) QC Batch: 63473 Prep Batch: 54173		nalyzed: 2009-09-03 paration: 2009-09-03		Analyzed By: MD Prepared By: MD	
Parameter	Flag	Result	Units	$\begin{array}{c} \text{Reporting} \\ \text{Limits} \end{array}$	
Total Dissolved Solids		< 5.00	m mg/L	5	
Method Blank (1)					
QC Batch: 63485 Prep Batch: 54188		nalyzed: 2009-09-11 eparation: 2009-09-11		Analyzed By: KV Prepared By: KV	

 ${\bf Result}$

< 0.0350

 Units

mg/L

Flag

Parameter

Nitrate and Nitrite as N

Reporting

Limits

0.035

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HELSTF Diesel Spill Groundwater

Method Blank (1)						
QC Batch: 63527 Prep Batch: 54231		Date Analyzed: 2009-09-10 QC Preparation: 2009-09-10		Analyzed By: JG Prepared By: JG		
Parameter	Flag	Result	${ m Units}$	$\begin{array}{c} \text{Reporting} \\ \text{Limits} \end{array}$		
Hydroxide Alkalinity Carbonate Alkalinity Bicarbonate Alkalinity Total Alkalinity		<1.00 <1.00 <4.00 <4.00	mg/L as CaCo3 mg/L as CaCo3 mg/L as CaCo3 mg/L as CaCo3	$egin{array}{cccccccccccccccccccccccccccccccccccc$		
Method Blank (1)						
QC Batch: 63668 Prep Batch: 54356		Date Analyzed: 2009-09-01 QC Preparation: 2009-09-01		Analyzed By: JR Prepared By: JR		
Parameter	Flag	Result	Units	$rac{ m Reporting}{ m Limits}$		
Chloride		< 0.640	m mg/L	0.6404		
Method Blank (1)						
QC Batch: 63668 Prep Batch: 54356		Date Analyzed: 2009-09-01 QC Preparation: 2009-09-01		Analyzed By: JR Prepared By: JR		
Parameter Sulfate	Flag	Result <0.504	Units mg/L	Reporting Limits 0.5038		
Durace		(0.501	1116/12	0.0000		
Method Blank (1)						
QC Batch: 63674 Prep Batch: 54363		Date Analyzed: 2009-09-16 QC Preparation: 2009-09-16		Analyzed By: JR Prepared By: JR		
				Reporting		
Parameter Bromide	Flag	Result < 0.0394	$\frac{\rm Units}{\rm mg/L}$	Limits 0.0394		

Date Analyzed:

QC Preparation: 2009-09-16

2009-09-16

Analyzed By: JR

Prepared By: JR

QC Batch:

Prep Batch: 54363

63674

${\it Work~Order:~9090211} \\ {\it HELSTF~Diesel~Spill~Groundwater}$

				Reporting
Parameter	Flag	Result	Units	Limits
Fluoride		< 0.0434	m mg/L	0.0434

Method Blank (1)

QC Batch: 63699 Prep Batch: 54391 Date Analyzed: 2009-09-17 QC Preparation: 2009-09-17 Analyzed By: JR Prepared By: MD

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				Reporting
Parameter	Flag	Result	Units	Limits
Oil and Grease		< 3.60	m mg/L	3.6

Duplicate (1) Duplicated Sample: 208955

QC Batch: 63339 Prep Batch: 54063 Date Analyzed: 2009-09-01 QC Preparation: 2009-09-01 Analyzed By: JG Prepared By: JG

	Duplicate	\mathbf{Sample}				RPD
Param	Result	Result	Units	Dilution	RPD	Limit
pН	5.56	5.58	s.u.	1	0	1.1

Duplicate (1) Duplicated Sample: 208953

QC Batch: 63473 Prep Batch: 54173 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03

Analyzed By: MD Prepared By: MD

	Duplicate	\mathbf{Sample}				RPD
Param	Result	Result	Units	Dilution	RPD	Limit
Total Dissolved Solids	28800	29500	$_{ m mg/L}$	1	2	10

Duplicate (1) Duplicated Sample: 208953

QC Batch: 63527 Prep Batch: 54231 Date Analyzed: 2009-09-10 QC Preparation: 2009-09-10

Analyzed By: JG Prepared By: JG

	Duplicate	\mathbf{Sample}				RPD
Param	Result	Result	Units	$\operatorname{Dilution}$	RPD	Limit
Hydroxide Alkalinity	<1.00	<1.00	mg/L as CaCo3	1	0	20
Carbonate Alkalinity	< 1.00	< 1.00	mg/L as $CaCo3$	1	0	20
Bicarbonate Alkalinity	262	260	mg/L as CaCo3	1	1	20
Total Alkalinity	262	260	mg/L as $CaCo3$	1	1	20

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Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Silver	0.124	${ m mg/L}$	1	0.125	< 0.00111	99	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit	RPD	Limit
Total Silver	0.124	mg/L	1	0.125	< 0.00111	99	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Total Aluminum	0.953	$_{ m mg/L}$	1	1.00	< 0.00301	95	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Aluminum	0.951	mg/L	1	1.00	< 0.00301	95	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Arsenic	0.478	$\mathrm{mg/L}$	1	0.500	< 0.00448	96	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Arsenic	0.476	$\mathrm{mg/L}$	1	0.500	< 0.00448	95	85 - 115	0	20

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

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	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	$\mathrm{Dil}.$	${f Amount}$	Result	Rec .	Limit
Total Barium	1.03	$\mathrm{mg/L}$	1	1.00	< 0.00105	103	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		Rec .		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Barium	1.03	$\mathrm{mg/L}$	1	1.00	< 0.00105	103	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Total Beryllium	0.0248	$_{ m mg/L}$	1	0.0250	< 0.000450	99	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil .	$\mathbf{A}\mathbf{mount}$	Result	Rec.	Limit	RPD	Limit
Total Beryllium	0.0256	mg/L	1	0.0250	< 0.000450	102	85 - 115	3	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	${f Limit}$
Total Cadmium	0.250	$\mathrm{mg/L}$	1	0.250	< 0.000303	100	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Cadmium	0.250	mg/L	1	0.250	< 0.000303	100	85 - 115	0	20

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Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Cobalt	0.246	${ m mg/L}$	1	0.250	< 0.000822	98	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Cobalt	0.246	mg/L	1	0.250	< 0.000822	98	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Total Chromium	0.0985	$_{ m mg/L}$	1	0.100	< 0.000583	98	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		Rec.		RPD
Param	Result	Units	Dil.	A mount	Result	Rec.	Limit	RPD	Limit
Total Chromium	0.0980	mg/L	1	0.100	< 0.000583	98	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	${f Limit}$
Total Copper	0.131	m mg/L	1	0.125	< 0.000843	105	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		Rec.		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Copper	0.133	mg/L	1	0.125	< 0.000843	106	85 - 115	2	20

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Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${ m Amount}$	Result	$\mathrm{Rec}.$	Limit
Total Iron	0.570	${ m mg/L}$	1	0.500	< 0.000872	114	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		Rec .		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
Total Iron	0.516	mg/L	1	0.500	< 0.000872	103	85 - 115	10	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Total Manganese	0.253	$_{ m mg/L}$	1	0.250	< 0.000305	101	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Manganese	0.254	mg/L	1	0.250	< 0.000305	102	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	${f Limit}$
Total Molybdenum	0.544	$\mathrm{mg/L}$	1	0.500	< 0.00119	109	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Molybdenum	0.543	mg/L	1	0.500	< 0.00119	109	85 - 115	0	20

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Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

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	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	$\mathrm{Rec}.$	Limit
Total Nickel	0.252	${ m mg/L}$	1	0.250	< 0.00121	101	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Nickel	0.251	mg/L	1	0.250	< 0.00121	100	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Total Phosphorous	0.473	$_{ m mg/L}$	1	0.500	< 0.00289	95	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Phosphorous	0.470	mg/L	1	0.500	< 0.00289	94	85 - 115	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	$\mathrm{Rec}.$	${f Limit}$
Total Lead	0.468	$\mathrm{mg/L}$	1	0.500	< 0.00326	94	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Lead	0.464	mg/L	1	0.500	< 0.00326	93	85 - 115	1	20

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

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	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Total Antimony	0.249	${ m mg/L}$	1	0.250	< 0.00440	100	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Antimony	0.248	mg/L	1	0.250	< 0.00440	99	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	$\mathrm{Dil}.$	${f Amount}$	Result	Rec .	Limit
Total Selenium	0.442	$_{ m mg/L}$	1	0.500	< 0.00508	88	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Selenium	0.443	mg/L	1	0.500	< 0.00508	89	85 - 115	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	${f Limit}$
Total Thallium	0.503	$\mathrm{mg/L}$	1	0.500	< 0.00488	101	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Thallium	0.503	mg/L	1	0.500	< 0.00488	101	85 - 115	0	20

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

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	LCS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Vanadium	0.246	$\mathrm{mg/L}$	1	0.250	< 0.000426	98	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Vanadium	0.248	mg/L	1	0.250	< 0.000426	99	85 - 115	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	$_{ m LCS}$			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Total Zinc	0.251	$_{ m mg/L}$	1	0.250	< 0.000465	100	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	$_{ m LCSD}$			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil .	$\mathbf{A}\mathbf{mount}$	Result	Rec.	Limit	RPD	Limit
Total Zinc	0.253	mg/L	1	0.250	< 0.000465	101	85 - 115	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63227 Prep Batch: 53960 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: TP Prepared By: TP

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	$\mathrm{Rec}.$	Limit
Total Mercury	0.000960	m mg/L	1	0.00100	< 0.0000329	96	90.3 - 108

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		Rec .		RPD
Param	Result	Units	Dil.	Amount	Result	Rec .	Limit	RPD	Limit
Total Mercury	0.000960	mg/L	1	0.00100	< 0.0000329	96	90.3 - 108	0	20

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Laboratory Control Spike (LCS-1)

QC Batch: 63239 Date Analyzed: 2009-09-03 Analyzed By: ER
Prep Batch: 53976 QC Preparation: 2009-09-03 Prepared By: ER

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	$\mathrm{Dil}.$	${f Amount}$	Result	Rec .	Limit
GRO	1.08	mg/L	1	1.00	< 0.152	108	75.5 - 118

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit	RPD	Limit
GRO	1.05	$_{ m mg/L}$	1	1.00	< 0.152	105	75.5 - 118	3	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCS	LCSD			Spike	LCS	LCSD	$\mathrm{Rec.}$
Surrogate	Result	Result	Units	Dil.	${f Amount}$	Rec .	Rec .	Limit
Trifluorotoluene (TFT)	0.109	0.104	$\mathrm{mg/L}$	1	0.100	109	104	78.2 - 121
4-Bromofluorobenzene (4-BFB)	0.106	0.104	$_{ m mg/L}$	1	0.100	106	104	82.2 - 118

Laboratory Control Spike (LCS-1)

QC Batch: 63261 Date Analyzed: 2009-09-04 Analyzed By: MN
Prep Batch: 53996 QC Preparation: 2009-09-02 Prepared By: MN

		LCS			Spike	Matrix		Rec .
Param		Result	Units	Dil.	${f Amount}$	Result	$\mathrm{Rec}.$	Limit
Phenol		0.0205	$\mathrm{mg/L}$	1	0.0800	< 0.000509	26	10 - 66.5
2-Chlorophenol		0.0411	$\mathrm{mg/L}$	1	0.0800	< 0.000537	51	11.2 - 108
1,4-Dichlorobenzene (para)		0.0378	$\mathrm{mg/L}$	1	0.0800	< 0.000440	47	16 - 101
N-Nitrosodi-n-propylamine		0.0474	$\mathrm{mg/L}$	1	0.0800	< 0.000732	59	10 - 142
1,2,4-Trichlorobenzene		0.0389	$\mathrm{mg/L}$	1	0.0800	< 0.000404	49	18 - 118
Naphthalene		0.0410	$\mathrm{mg/L}$	1	0.0800	< 0.000489	51	20.2 - 114
4-Chloro-3-methylphenol		0.0507	$\mathrm{mg/L}$	1	0.0800	< 0.000522	63	21.5 - 125
${ m Acenaphthylene}$		0.0474	$\mathrm{mg/L}$	1	0.0800	< 0.000586	59	25.8 - 121
${ m Acenaphthene}$		0.0474	$\mathrm{mg/L}$	1	0.0800	< 0.000423	59	33.5 - 122
4-Nitrophenol		0.0191	$\mathrm{mg/L}$	1	0.0800	< 0.00185	24	10 - 125
2,4-Dinitrotoluene		0.0543	$\mathrm{mg/L}$	1	0.0800	< 0.000911	68	53 - 130
Fluorene		0.0494	$\mathrm{mg/L}$	1	0.0800	< 0.000648	62	44.6 - 117
${ m Pentachlorophenol}$		0.0373	$\mathrm{mg/L}$	1	0.0800	< 0.000435	47	10 - 139
${ m Anthracene}$		0.0511	$\mathrm{mg/L}$	1	0.0800	< 0.000428	64	57.5 - 115
Phenanthrene		0.0509	$\mathrm{mg/L}$	1	0.0800	< 0.000548	64	55.5 - 118
Fluoranthene		0.0545	$\mathrm{mg/L}$	1	0.0800	< 0.000632	68	57 - 122
Pyrene		0.0468	$\mathrm{mg/L}$	1	0.0800	< 0.000723	58	58.5 - 130
$\operatorname{Benzo}(\operatorname{a})\operatorname{anthracene}$	2	0.0463	m mg/L	1	0.0800	< 0.000527	58	63.4 - 109
Chrysene		0.0508	m mg/L	1	0.0800	< 0.000638	64	54.7 - 114

 $continued \dots$

²Spike analyte out of control limits. Results biased low. •

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control spikes continued

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		LCS			Spike	Matrix		$\mathrm{Rec}.$
Param		Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Benzo(b)fluoranthene	3	0.0477	m mg/L	1	0.0800	< 0.000879	60	64.8 - 120
Benzo(k)fluoranthene		0.0619	m mg/L	1	0.0800	< 0.000845	77	70.3 - 114
Benzo(a)pyrene		0.0575	${ m mg/L}$	1	0.0800	< 0.00167	72	63.7 - 120
Indeno(1,2,3-cd)pyrene		0.0568	${ m mg/L}$	1	0.0800	< 0.000862	71	65.4 - 119
${ m Dibenzo}({ m a,h}) { m anthracene}$		0.0572	m mg/L	1	0.0800	< 0.000809	72	68.7 - 117
$\operatorname{Benzo}(g,h,i)$ perylene		0.0598	$\mathrm{mg/L}$	1	0.0800	< 0.000949	75	57.2 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Phenol	0.0210	mg/L	1	0.0800	< 0.000509	26	10 - 66.5	2	20
2-Chlorophenol	0.0432	$\mathrm{mg/L}$	1	0.0800	< 0.000537	54	11.2 - 108	5	20
1,4-Dichlorobenzene (para)	0.0398	$\mathrm{mg/L}$	1	0.0800	< 0.000440	50	16 - 101	5	20
N-Nitrosodi-n-propylamine	0.0506	$\mathrm{mg/L}$	1	0.0800	< 0.000732	63	10 - 142	6	20
1,2,4-Trichlorobenzene	0.0413	$\mathrm{mg/L}$	1	0.0800	< 0.000404	52	18 - 118	6	20
Naphthalene	0.0436	$\mathrm{mg/L}$	1	0.0800	< 0.000489	54	20.2 - 114	6	20
4-Chloro-3-methylphenol	0.0534	$\mathrm{mg/L}$	1	0.0800	< 0.000522	67	21.5 - 125	5	20
${ m Acenapht}$ hylene	0.0515	$\mathrm{mg/L}$	1	0.0800	< 0.000586	64	25.8 - 121	8	20
${\it Acenaphthene}$	0.0505	$\mathrm{mg/L}$	1	0.0800	< 0.000423	63	33.5 - 122	6	20
4-Nitrophenol	0.0220	$\mathrm{mg/L}$	1	0.0800	< 0.00185	28	10 - 125	14	20
2,4-Dinitrotoluene	0.0577	$\mathrm{mg/L}$	1	0.0800	< 0.000911	72	53 - 130	6	20
Fluorene	0.0533	$\mathrm{mg/L}$	1	0.0800	< 0.000648	67	44.6 - 117	8	20
${ m Pentachlorophenol}$	0.0388	$\mathrm{mg/L}$	1	0.0800	< 0.000435	48	10 - 139	4	20
${ m Anthracene}$	0.0530	$\mathrm{mg/L}$	1	0.0800	< 0.000428	66	57.5 - 115	4	20
Phenanthrene	0.0521	$\mathrm{mg/L}$	1	0.0800	< 0.000548	65	55.5 - 118	2	20
Fluoranthene	0.0564	$\mathrm{mg/L}$	1	0.0800	< 0.000632	70	57 - 122	3	20
Pyrene	0.0495	$\mathrm{mg/L}$	1	0.0800	< 0.000723	62	58.5 - 130	6	20
Benzo(a)anthracene 4	0.0486	$\mathrm{mg/L}$	1	0.0800	< 0.000527	61	63.4 - 109	5	20
Chrysene	0.0541	$\mathrm{mg/L}$	1	0.0800	< 0.000638	68	54.7 - 114	6	20
$\operatorname{Benzo}(\operatorname{b})$ fluoranthene	0.0533	$\mathrm{mg/L}$	1	0.0800	< 0.000879	67	64.8 - 120	11	20
$\operatorname{Benzo}(k)$ fluoranthene	0.0696	$\mathrm{mg/L}$	1	0.0800	< 0.000845	87	70.3 - 114	12	20
$\mathrm{Benzo}(\mathrm{a})\mathrm{pyrene}$	0.0630	$\mathrm{mg/L}$	1	0.0800	< 0.00167	79	63.7 - 120	9	20
$\operatorname{Indeno}(1,2,3\text{-cd})$ pyrene	0.0615	$\mathrm{mg/L}$	1	0.0800	< 0.000862	77	65.4 - 119	8	20
${ m Dibenzo}({ m a,h}) { m anthracene}$	0.0630	$\mathrm{mg/L}$	1	0.0800	< 0.000809	79	68.7 - 117	10	20
$\mathrm{Benzo}(\mathrm{g,h,i})$ perylene	0.0643	$\mathrm{mg/L}$	1	0.0800	< 0.000949	80	57.2 - 125	7	20

Surrogate	$egin{array}{c} ext{LCS} \ ext{Result} \end{array}$	$egin{array}{c} ext{LCSD} \ ext{Result} \end{array}$	Units	Dil.	$rac{ ext{Spike}}{ ext{Amount}}$	$\frac{\text{LCS}}{\text{Rec.}}$	$\begin{array}{c} { m LCSD} \\ { m Rec.} \end{array}$	$egin{array}{c} { m Rec.} \\ { m Limit} \end{array}$
2-Fluorophenol	0.0260	0.0278	mg/L	1	0.0800	32	35	10 - 53.1
Phenol-d5	0.0198	0.0210	mg/L	1	0.0800	25	26	10 - 36.9
${ m Nitrobenzene-d5}$	0.0419	0.0439	mg/L	1	0.0800	52	55	23.8 - 108
2-Fluorobiphenyl	0.0429	0.0466	mg/L	1	0.0800	54	58	15.9 - 127
2,4,6-Tribromophenol	0.0565	0.0609	mg/L	1	0.0800	71	76	10 - 123
Terphenyl-d14	0.0461	0.0494	m mg/L	1	0.0800	58	62	17.2 - 160

 $^{^3}$ Spike analyte out of control limits. Results biased low. \bullet 4 Spike analyte out of control limits. Results biased low. \bullet

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Laboratory Control Spike (LCS-1)

QC Batch: 63307 Prep Batch: 54035 Date Analyzed: 2009-09-07 QC Preparation: 2009-09-04 Analyzed By: Prepared By:

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	$\mathrm{Rec}.$	Limit
DRO	27.0	${ m mg/L}$	1	25.0	< 0.876	108	78.6 - 154

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit	RPD	Limit
DRO	28.3	$\mathrm{mg/L}$	1	25.0	< 0.876	113	78.6 - 154	5	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCS	LCSD			Spike	LCS	LCSD	$\mathrm{Rec}.$
Surrogate	Result	Result	Units	Dil.	${f Amount}$	$\mathrm{Rec.}$	Rec .	Limit
n-Triacontane	8.66	8.90	mg/L	1	10.0	87	89	57.3 - 151

Laboratory Control Spike (LCS-1)

QC Batch: 63327 Prep Batch: 54055 Date Analyzed: 2009-09-01 QC Preparation: 2009-09-01 Analyzed By: MD Prepared By: MD

	$_{ m LCS}$			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Hexavalent Chromium	0.489	mg/L	1	0.500	< 0.00594	98	95.4 - 105

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec.}$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Hexavalent Chromium	0.497	mg/L	1	0.500	< 0.00594	99	95.4 - 105	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63340 Prep Batch: 54065 Date Analyzed: 2009-09-08 QC Preparation: 2009-09-08 Analyzed By: KV Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Organic Carbon	48.4	${ m mg/L}$	1	50.0	< 0.401	97	89.5 - 114

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

 $continued \dots$

control spikes continued									
-	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
	$_{ m LCSD}$			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	$\mathrm{Rec}.$	Limit	RPD	Limit
Total Organic Carbon	49.4	mg/L	1	50.0	< 0.401	99	89.5 - 114	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63355 Prep Batch: 53951 Date Analyzed: 2009-09-09 QC Preparation: 2009-09-03

Analyzed By: RR Prepared By: KV

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	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Total Calcium	54.2	mg/L	1	50.0	< 0.117	108	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Calcium	50.5	$\mathrm{mg/L}$	1	50.0	< 0.117	101	85 - 115	7	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63355 Prep Batch: 53951 Date Analyzed: 2009-09-09 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Total Potassium	53.3	m mg/L	1	50.0	< 0.172	107	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		Rec.		RPD
Param	Result	Units	Dil .	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Potassium	50.2	mg/L	1	50.0	< 0.172	100	85 - 115	6	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63355 Prep Batch: 53951 Date Analyzed: 2009-09-09 QC Preparation: 2009-09-03

Analyzed By: RR Prepared By: KV

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	$_{ m LCS}$			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Magnesium	52.2	mg/L	1	50.0	< 0.160	104	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Magnesium	50.1	$\mathrm{mg/L}$	1	50.0	< 0.160	100	85 - 115	4	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Total Sodium	53.2	mg/L	1	50.0	< 0.0500	106	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Sodium	50.6	mg/L	1	50.0	< 0.0500	101	85 - 115	5	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63425 Date Analyzed: 2009-09-10 Analyzed By: DS
Prep Batch: 54137 QC Preparation: 2009-09-04 Prepared By: DS

Param	$rac{ ext{LCS}}{ ext{Result}}$	Units	Dil.	${ m Spike} \ { m Amount}$	$rac{ m Matrix}{ m Result}$	${ m Rec.}$	$egin{array}{c} { m Rec.} \\ { m Limit} \end{array}$
HMX	2.41	$\mu \mathrm{g/L}$	1	2.50	< 0.123	96	63.5 - 125
RDX	2.37	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.298	95	74.5 - 124
1,3,5-Trinitrobenzene	2.38	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.339	95	54.1 - 131
1,3-Dinitrobenzene	2.45	$\mu { m g}/{ m L}$	1	2.50	< 0.389	98	72 - 112
Nitrobenzene	2.50	$\mu { m g}/{ m L}$	1	2.50	< 0.379	100	72.5 - 126
Tetryl	2.31	$\mu { m g}/{ m L}$	1	2.50	< 0.413	92	35.9 - 149
TNT	2.37	$\mu { m g}/{ m L}$	1	2.50	< 0.464	95	40.7 - 129
4-Amino-DNT	2.31	$\mu { m g}/{ m L}$	1	2.50	< 0.319	92	80 - 120
2-Amino-DNT	2.51	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.391	100	80 - 120
2,6-DNT	2.19	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.323	88	80 - 120
2,4-DNT	2.52	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.366	101	80 - 120
2-NT	2.47	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.379	99	49.8 - 139
4-NT	2.07	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.398	83	56.3 - 141
3-NT	2.37	$\mu { m g}/{ m L}$	1	2.50	< 0.346	95	66.2 - 129

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KV

	LCSD			Spike	Matrix		Rec.		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
HMX	2.43	$\mu \mathrm{g/L}$	1	2.50	< 0.123	97	63.5 - 125	1	20
RDX	2.33	$\mu { m g}/{ m L}$	1	2.50	< 0.298	93	74.5 - 124	2	20
1,3,5-Trinitrobenzene	2.38	$\mu { m g}/{ m L}$	1	2.50	< 0.339	95	54.1 - 131	0	20
1,3-Dinitrobenzene	2.41	$\mu { m g}/{ m L}$	1	2.50	< 0.389	96	72 - 112	2	20
Nitrobenzene	2.46	$\mu { m g}/{ m L}$	1	2.50	< 0.379	98	72.5 - 126	2	20
Tetryl	2.26	$\mu { m g}/{ m L}$	1	2.50	< 0.413	90	35.9 - 149	2	20
TNT	2.33	$\mu { m g}/{ m L}$	1	2.50	< 0.464	93	40.7 - 129	2	20
4-Amino-DNT	2.37	$\mu { m g}/{ m L}$	1	2.50	< 0.319	95	80 - 120	3	20
2-Amino-DNT	2.64	$\mu { m g}/{ m L}$	1	2.50	< 0.391	106	80 - 120	5	20
2,6-DNT	2.29	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.323	92	80 - 120	4	20
2,4-DNT	2.60	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.366	104	80 - 120	3	20
2-NT	2.34	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.379	94	49.8 - 139	5	20
4-NT	2.26	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.398	90	56.3 - 141	9	20
3-NT	2.43	$\mu \mathrm{g}/\mathrm{L}$	1	2.50	< 0.346	97	66.2 - 129	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCS	LCSD			Spike	LCS	LCSD	$\mathrm{Rec}.$
$\mathbf{Surrogate}$	Result	Result	Units	Dil.	${f Amount}$	Rec .	$\mathrm{Rec}.$	Limit
1,2-Dinitrobenzene	2.29	2.15	$\mu { m g/L}$	1	2.50	92	86	53 - 134

Laboratory Control Spike (LCS-1)

QC Batch: Date Analyzed: Analyzed By: RR 63461 2009-09-11 Prep Batch: 54153 QC Preparation: 2009-09-11 Prepared By:

	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Dissolved Chromium	0.0990	$_{ m mg/L}$	1	0.100	< 0.000583	99	85 - 115

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Dissolved Chromium	0.0950	mg/L	1	0.100	< 0.000583	95	85 - 115	4	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63668 Date Analyzed: Analyzed By: JR 2009-09-01 Prep Batch: 54356 QC Preparation: 2009-09-01 Prepared By: JR

	$_{ m LCS}$			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Chloride	23.1	$_{ m mg/L}$	1	25.0	< 0.640	92	90 - 110

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	LCSD			Spike	Matrix		Rec.		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
Chloride	23.0	mg/L	1	25.0	< 0.640	92	90 - 110	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63668 Date Analyzed: 2009-09-01 Analyzed By: JR
Prep Batch: 54356 QC Preparation: 2009-09-01 Prepared By: JR

LCS Spike Rec. Matrix Param Result Units Dil. Amount Result Rec. Limit Sulfate 22.8 25.0 < 0.50491 90 - 110 mg/L1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil .	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Sulfate	22.6	mg/L	1	25.0	< 0.504	90	90 - 110	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63674 Date Analyzed: 2009-09-16 Analyzed By: JR
Prep Batch: 54363 QC Preparation: 2009-09-16 Prepared By: JR

LCS Spike MatrixRec. Param Result Units Dil. Amount Result Rec. Limit Bromide 4.82 mg/L5.00 < 0.039496 90 - 110

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		Rec .		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Bromide	4.81	mg/L	1	5.00	< 0.0394	96	90 - 110	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

	LCS			Spike	Matrix		$\mathrm{Rec.}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Fluoride	4.90	$\mathrm{mg/L}$	1	5.00	< 0.0434	98	90 - 110

	I CCD			G :1	3.6		D		DDD
	LCSD			Spike	Matrix		$\mathrm{Rec}.$		$_{ m RPD}$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Fluoride	4.89	mg/L	1	5.00	< 0.0434	98	90 - 110	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Laboratory Control Spike (LCS-1)

QC Batch: 63699 Prep Batch: 54391 Date Analyzed: 2009-09-17 QC Preparation: 2009-09-17 Analyzed By: JR Prepared By: MD

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	LCS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${\bf Amount}$	Result	Rec .	Limit
Oil and Grease	39.5	$\mathrm{mg/L}$	1	40.0	< 3.60	99	78 - 114

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	LCSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Oil and Grease	35.2	${ m mg/L}$	1	40.0	< 3.60	88	78 - 114	12	18

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Prep Batch: 53951 Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	$\mathrm{Dil}.$	${f Amount}$	Result	Rec .	Limit
Total Silver	0.137	m mg/L	1	0.125	< 0.00111	110	75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
Total Silver	0.135	mg/L	1	0.125	< 0.00111	108	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

 QC Batch:
 63221
 Da

 Prep Batch:
 53951
 QC

Date Analyzed: 2009-09-03 QC Preparation: 2009-09-03 Analyzed By: RR Prepared By: KV

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Aluminum	1.26	${ m mg/L}$	1	1.00	0.065	120	75 - 125

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	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Aluminum	1.24	mg/L	1	1.00	0.065	118	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSMatrix Spike Rec. Param Result Units Dil. Amount Result Rec. Limit Total Arsenic 0.506 0.500 < 0.0044875 - 125 mg/L101 1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Arsenic	0.496	mg/L	1	0.500	< 0.00448	99	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike MatrixRec. Param Result Units Dil. Amount Result Rec. Limit 0.008 Total Barium 1.00 mg/L1.00 99 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Barium	0.988	$_{ m mg/L}$	1	1.00	0.008	98	75 - 125	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit Total Beryllium 0.0250 < 0.000450 102 0.0254mg/L75 - 125

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	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Beryllium	0.0247	mg/L	1	0.0250	< 0.000450	99	75 - 125	3	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit Total Cadmium 0.227 0.250 < 0.00030375 - 125 mg/L91 1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Cadmium	0.222	mg/L	1	0.250	< 0.000303	89	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit Total Cobalt 0.230 mg/L0.250< 0.00082292 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Cobalt	0.226	mg/L	1	0.250	< 0.000822	90	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit 0.100 Total Chromium 0.111mg/L0.01596 75 - 125

${\it Work~Order:~9090211} \\ {\it HELSTF~Diesel~Spill~Groundwater}$

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'									
	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil .	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Chromium	0.108	mg/L	1	0.100	0.015	93	75 - 125	3	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit Total Copper 0.1440.125 < 0.00084375 - 125 mg/L115 1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Copper	0.140	mg/L	1	0.125	< 0.000843	112	75 - 125	3	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit Total Iron 0.488 mg/L0.500< 0.00087298 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Iron	0.462	mg/L	1	0.500	< 0.000872	92	75 - 125	6	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit 0.2350.250 < 0.000305 Total Manganese mg/L94 75 - 125

Total Molybdenum

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	MSD			Spike	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Manganese	0.231	mg/L	1	0.250	< 0.000305	92	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

mg/L

MS Spike Matrix Rec.
Param Result Units Dil. Amount Result Rec. Limit

1

0.500

0.211

112

75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

0.771

MSD Spike Matrix RPD Rec. Param Result Units Dil. Amount Result Rec. Limit RPD Limit Total Molybdenum 109 20 0.758mg/L0.5000.211 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Dil. Result Units A mount Result Rec. Limit Total Nickel 0.230 mg/L0.250< 0.0012192 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSD RPD Spike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit RPD Limit 0.225 0.250 20 Total Nickel mg/L< 0.00121 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Result Units Dil. Amount Result Rec. Limit 0.5000.493Total Phosphorous mg/L0.00498 75 - 125

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	MOD			G 11	3.5		D.		DDD
	MSD			${ m Spike}$	Matrix		${ m Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit	RPD	Limit
Total Phosphorous	0.485	mg/L	1	0.500	0.004	96	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSMatrix Spike Rec. Param Result Units Dil. Amount Result Rec. Limit Total Lead 0.412 0.500< 0.0032675 - 125 mg/L82 1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSD Spike Matrix RPD Rec. Param Result Units Dil. A mount Result Rec. Limit RPD Limit Total Lead 0.403 20 mg/L0.500< 0.0032675 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Dil. Result Units A mount Result Rec. Limit Total Antimony 0.246mg/L0.250< 0.0044098 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSD RPD Spike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit RPD Limit 0.242 0.25020 Total Antimony mg/L< 0.00440 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit 0.502< 0.00508100 Total Selenium mg/L0.50075 - 125

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KV

	MSD			Spike	Matrix		$\mathrm{Rec.}$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Selenium	0.487	mg/L	1	0.500	< 0.00508	97	75 - 125	3	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

2009-09-03 QC Batch: 63221 Date Analyzed: Analyzed By: RR Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By:

MSMatrix Spike Rec. Param Result Units Dil. Amount Result Rec. Limit Total Thallium 0.4390.500 < 0.0048875 - 125 mg/L88 1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Thallium	0.437	mg/L	1	0.500	< 0.00488	87	75 - 125	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: Date Analyzed: 2009-09-03 Analyzed By: RR 63221 Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike MatrixRec. Param Units Result Dil. Amount Result Rec. Limit Total Vanadium 0.265mg/L0.2500.01799 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	$_{ m Limit}$	RPD	Limit
Total Vanadium	0.260	$_{ m mg/L}$	1	0.250	0.017	97	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR Prep Batch: 2009-09-03 Prepared By: KV53951QC Preparation:

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit 0.2580.250 < 0.000465 103 Total Zinc mg/L75 - 125

	MSD			Spike	Matrix		Rec.		RPD
Param	Result	Units	Dil.	Amount	Result	Rec.	m Limit	RPD	Limit
Total Zinc	0.253	mg/L	1	0.250	< 0.000465	101	75 - 125	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 209000

2009-09-03 QC Batch: 63227Date Analyzed: Prep Batch: 53960

Analyzed By: TP QC Preparation: 2009-09-03 Prepared By: TP

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		MS			Spike	Matrix		$\mathrm{Rec}.$
Param		Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Mercury	5	0.000850	mg/L	1	0.00100	6e-05	79	80 - 116

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec.}$		RPD
Param	Result	Units	Dil.	Amount	Result	$\mathrm{Rec}.$	Limit	RPD	Limit
Total Mercury	0.000860	mg/L	1	0.00100	6e-05	80	80 - 116	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 209098

QC Batch: Date Analyzed: 2009-09-03 Analyzed By: ER 63239 Prep Batch: 53976QC Preparation: 2009-09-03 Prepared By:

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	$\mathrm{Rec.}$	Limit
GRO	1.11	$_{ m mg/L}$	1	1.00	< 0.152	111	48.4 - 136

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

		MSD			Spike	Matrix		Rec .		RPD
Param		Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
GRO	6	0.768	$_{ m mg/L}$	1	1.00	< 0.152	77	48.4 - 136	36	20

		MS	MSD			Spike	MS	MSD	$\mathrm{Rec.}$
$\operatorname{Surrogate}$		Result	Result	Units	Dil.	${f Amount}$	$\mathrm{Rec.}$	Rec .	Limit
Trifluorotoluene (TFT)	7	0.112	0.0584	$\mathrm{mg/L}$	1	0.1	112	58	70.3 - 129
4-Bromofluorobenzene (4-BFB)	8	0.111	0.0576	mg/L	1	0.1	111	58	82.5 - 118

⁵Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

⁶MS/MSD RPD out of RPD Limits. Use LCS/LCSD to demonstrate analysis is under control.

⁷ Matrix spike recovery out of control limits. Use LCS/LCSD to demonstrate analysis is under control.

⁸ Matrix spike recovery out of control limits. Use LCS/LCSD to demonstrate analysis is under control.

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Spiked Sample: 208408 Matrix Spike (MS-1)

2009-09-04 QC Batch: 63261Date Analyzed: Analyzed By: MN Prep Batch: 53996QC Preparation: 2009-09-02Prepared By: MN

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Phenol	0.0194	$\mathrm{mg/L}$	1.02	0.0800	< 0.000519	24	10 - 66.5
2-Chlorophenol	0.0437	$\mathrm{mg/L}$	1.02	0.0800	< 0.000548	55	11.2 - 108
1,4-Dichlorobenzene (para)	0.0354	$\mathrm{mg/L}$	1.02	0.0800	< 0.000449	44	16 - 101
N-Nitrosodi-n-propylamine	0.0490	$\mathrm{mg/L}$	1.02	0.0800	< 0.000747	61	10 - 142
1,2,4-Trichlorobenzene	0.0389	$\mathrm{mg/L}$	1.02	0.0800	< 0.000412	49	18 - 108
${f Naphthalene}$	0.0420	$\mathrm{mg/L}$	1.02	0.0800	< 0.000499	52	20.2 - 114
4-Chloro-3-methylphenol	0.0558	$\mathrm{mg/L}$	1.02	0.0800	< 0.000532	70	21.5 - 125
${ m Acenapht}$ hylene	0.0543	$\mathrm{mg/L}$	1.02	0.0800	< 0.000598	68	25.8 - 121
${ m Acenapht}$ hene	0.0538	$\mathrm{mg/L}$	1.02	0.0800	< 0.000431	67	33.5 - 122
4-Nitrophenol	0.0218	$\mathrm{mg/L}$	1.02	0.0800	< 0.00189	27	10 - 125
2,4-Dinitrotoluene	0.0604	$\mathrm{mg/L}$	1.02	0.0800	< 0.000929	76	53 - 130
Fluorene	0.0572	$\mathrm{mg/L}$	1.02	0.0800	< 0.000661	72	44.6 - 117
${ m Pentachlorophenol}$	0.0394	${ m mg/L}$	1.02	0.0800	< 0.000444	49	10 - 139
${ m Anthracene}$	0.0557	${ m mg/L}$	1.02	0.0800	< 0.000436	70	57.5 - 115
Phenanthrene	0.0558	${ m mg/L}$	1.02	0.0800	< 0.000559	70	55.5 - 118
Fluoranthene	0.0588	$\mathrm{mg/L}$	1.02	0.0800	< 0.000645	74	57 - 122
Pyrene	0.0524	$\mathrm{mg/L}$	1.02	0.0800	< 0.000737	66	58.5 - 130
$\operatorname{Benzo}(\operatorname{a}) \operatorname{anthracene}$	0.0508	$\mathrm{mg/L}$	1.02	0.0800	< 0.000538	64	63.4 - 109
Chrysene	0.0568	$\mathrm{mg/L}$	1.02	0.0800	< 0.000651	71	54.7 - 114
$\operatorname{Benzo}(\operatorname{b})$ fluoranthene	0.0546	${ m mg/L}$	1.02	0.0800	< 0.000896	68	64.8 - 120
$\operatorname{Benzo}(k)$ fluoranthene	0.0721	${ m mg/L}$	1.02	0.0800	< 0.000862	90	70.3 - 114
$\mathrm{Benzo}(\mathrm{a})\mathrm{pyrene}$	0.0675	$\mathrm{mg/L}$	1.02	0.0800	< 0.00170	84	63.7 - 120
Indeno(1,2,3-cd)pyrene	0.0643	$\mathrm{mg/L}$	1.02	0.0800	< 0.000879	80	65.4 - 119
${ m Dibenzo}({ m a,h}) { m anthracene}$	0.0634	$\mathrm{mg/L}$	1.02	0.0800	< 0.000825	79	68.7 - 117
Benzo(g,h,i)perylene	0.0664	$\mathrm{mg/L}$	1.02	0.0800	< 0.000968	83	57.2 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			$_{ m Spike}$	Matrix		$\mathrm{Rec.}$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit	RPD	Limit
Phenol	0.0182	$\mathrm{mg/L}$	1.02	0.0800	< 0.000519	23	10 - 66.5	6	20
2-Chlorophenol	0.0398	${ m mg/L}$	1.02	0.0800	< 0.000548	50	11.2 - 108	9	20
1,4-Dichlorobenzene (para)	0.0317	$\mathrm{mg/L}$	1.02	0.0800	< 0.000449	40	16 - 101	11	20
N-Nitrosodi-n-propylamine	0.0465	$\mathrm{mg/L}$	1.02	0.0800	< 0.000747	58	10 - 142	5	20
1,2,4-Trichlorobenzene	0.0344	${ m mg/L}$	1.02	0.0800	< 0.000412	43	18 - 108	12	20
Naphthalene	0.0363	$\mathrm{mg/L}$	1.02	0.0800	< 0.000499	45	20.2 - 114	15	20
4-Chloro-3-methylphenol	0.0496	$\mathrm{mg/L}$	1.02	0.0800	< 0.000532	62	21.5 - 125	12	20
${ m Acenapht}$ hylene	0.0479	$\mathrm{mg/L}$	1.02	0.0800	< 0.000598	60	25.8 - 121	12	20
${ m Acenaphthene}$	0.0468	$\mathrm{mg/L}$	1.02	0.0800	< 0.000431	58	33.5 - 122	14	20
4-Nitrophenol	0.0200	mg/L	1.02	0.0800	< 0.00189	25	10 - 125	9	20
2,4-Dinitrotoluene	0.0526	$\mathrm{mg/L}$	1.02	0.0800	< 0.000929	66	53 - 130	14	20
Fluorene	0.0497	$\mathrm{mg/L}$	1.02	0.0800	< 0.000661	62	44.6 - 117	14	20
Pentachlorophenol	0.0358	mg/L	1.02	0.0800	< 0.000444	45	10 - 139	10	20

 $continued \dots$

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matrix spikes continued . . .

		MSD			Spike	Matrix		$\mathrm{Rec.}$		RPD
Param		Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Anthracene		0.0503	mg/L	1.02	0.0800	< 0.000436	63	57.5 - 115	10	20
Phenanthrene		0.0492	$\mathrm{mg/L}$	1.02	0.0800	< 0.000559	62	55.5 - 118	13	20
Fluoranthene		0.0537	$\mathrm{mg/L}$	1.02	0.0800	< 0.000645	67	57 - 122	9	20
Pyrene	9	0.0457	$\mathrm{mg/L}$	1.02	0.0800	< 0.000737	57	58.5 - 130	14	20
$\operatorname{Benzo}(\operatorname{a})\operatorname{anthracene}$	10	0.0446	$\mathrm{mg/L}$	1.02	0.0800	< 0.000538	56	63.4 - 109	13	20
Chrysene		0.0485	$\mathrm{mg/L}$	1.02	0.0800	< 0.000651	61	54.7 - 114	16	20
${ m Benzo(b)}$ fluoranthene		0.0528	$\mathrm{mg/L}$	1.02	0.0800	< 0.000896	66	64.8 - 120	3	20
$\operatorname{Benzo}(k)$ fluoranthene	11	0.0573	$\mathrm{mg/L}$	1.02	0.0800	< 0.000862	72	70.3 - 114	23	20
$\mathrm{Benzo}(\mathrm{a})\mathrm{pyrene}$		0.0589	$\mathrm{mg/L}$	1.02	0.0800	< 0.00170	74	63.7 - 120	14	20
Indeno(1,2,3-cd)pyrene		0.0540	$\mathrm{mg/L}$	1.02	0.0800	< 0.000879	68	65.4 - 119	17	20
${ m Dibenzo(a,h)}$ anthracene	12	0.0536	$\mathrm{mg/L}$	1.02	0.0800	< 0.000825	67	68.7 - 117	17	20
$\mathrm{Benzo}(\mathrm{g,h,i})\mathrm{perylene}$		0.0557	$\mathrm{mg/L}$	1.02	0.0800	< 0.000968	70	57.2 - 125	18	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MS	MSD			Spike	MS	MSD	$\mathrm{Rec.}$
$\mathbf{Surrogate}$	Result	Result	Units	Dil.	${ m Amount}$	Rec .	Rec .	Limit
2-Fluorophenol	0.0262	0.0254	$\mathrm{mg/L}$	1.02	0.08	33	32	10 - 53.1
Phenol-d5	0.0194	0.0183	${ m mg/L}$	1.02	0.08	24	23	10 - 36.9
${ m Nitrobenzene-d5}$	0.0450	0.0403	${ m mg/L}$	1.02	0.08	56	50	23.8 - 108
2-Fluorobiphenyl	0.0445	0.0387	${ m mg/L}$	1.02	0.08	56	48	15.9 - 127
2,4,6-Tribromophenol	0.0653	0.0583	${ m mg/L}$	1.02	0.08	82	73	10 - 123
Terphenyl-d14	0.0522	0.0450	$\mathrm{mg/L}$	1.02	0.08	65	56	17.2 - 160

Matrix Spike (MS-1) Spiked Sample: 208666

QC Batch: Analyzed By: 63307 Date Analyzed: 2009-09-07 Prep Batch: 54035 QC Preparation: 2009-09-04 Prepared By:

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
DRO	24.3	m mg/L	1	25.0	< 0.876	97	54 - 144

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			$_{ m Spike}$	Matrix		Rec .		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
DRO	23.5	mg/L	1	25.0	< 0.876	94	54 - 144	3	20

	MS	MSD			Spike	MS	MSD	$\mathrm{Rec}.$
Surrogate	Result	Result	Units	Dil.	${f Amount}$	Rec.	$\mathrm{Rec.}$	Limit
n-Triacontane	10.9	10.9	m mg/L	1	10	109	109	57.3 - 151

⁹ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

¹⁰ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

¹¹MS/MSD RPD out of RPD Limits. Use LCS/LCSD to demonstrate analysis is under control.

¹² Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

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Matrix Spike (MS-1) Spiked Sample: 208955

QC Batch: 63327Date Analyzed: 2009-09-01 Analyzed By: MD Prep Batch: 54055QC Preparation: 2009-09-01 Prepared By: MD

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit
Hexavalent Chromium	0.561	${ m mg/L}$	1.11	0.556	< 0.00659	101	80.1 - 118

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		Rec .		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Hexavalent Chromium	0.573	mg/L	1.11	0.556	< 0.00659	103	80.1 - 118	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63340 Date Analyzed: 2009-09-08 Analyzed By: KV Prep Batch: 54065 Prepared By: QC Preparation: 2009-09-08 KV

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Organic Carbon	51.2	$_{ m mg/L}$	1	50.0	1.66	99	66.9 - 121

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec.}$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Organic Carbon	48.4	mg/L	1	50.0	1.66	93	66.9 - 121	6	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63355Date Analyzed: 2009-09-09 Analyzed By: RR Prep Batch: 53951QC Preparation: 2009-09-03 Prepared By: KV

	MS			Spike	Matrix		${ m Rec.}$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Total Calcium	574	mg/L	1	50.0	531	86	75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Calcium	576	mg/L	1	50.0	531	90	75 - 125	0	20

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RR

KV

Matrix Spike (MS-1) Spiked Sample: 208408

 QC Batch:
 63355
 Date Analyzed:
 2009-09-09
 Analyzed By:

 Prep Batch:
 53951
 QC Preparation:
 2009-09-03
 Prepared By:

MSSpike Matrix Rec. Param Result Units Dil. AmountResult Rec. Limit Total Potassium 150 99.9 100 75 - 125 mg/L50.0

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			$_{ m Spike}$	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Potassium	148	mg/L	1	50.0	99.9	96	75 - 125	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike MatrixRec. Dil. Limit Param Result Units AmountResult Rec. Total Magnesium 471 mg/L50.0425 92 75 - 125 1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${f Amount}$	Result	Rec.	Limit	RPD	Limit
Total Magnesium	472	mg/L	1	50.0	425	94	75 - 125	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RR
Prep Batch: 53951 QC Preparation: 2009-09-03 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Result Rec. Limit Amount Total Sodium 768 mg/L50.0 717102 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Sodium	776	mg/L	1	50.0	717	118	75 - 125	1	20

Work Order: 9090211

HELSTF Diesel Spill Groundwater

Matrix Spike (MS-1) Spiked Sample: 209099

QC Batch: 63370Date Analyzed: 2009-09-05 Prep Batch: 54092QC Preparation: 2009-09-05 Analyzed By: AH Prepared By: AH

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	MS			Spike	Matrix		Rec .
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
Ammonia-N	5.38	${ m mg/L}$	1	5.00	0.504	98	57.2 - 133

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Ammonia-N	5.26	mg/L	1	5.00	0.504	95	57.2 - 133	2	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 209099

QC Batch: 63391 Date Analyzed: 2009-09-07 Analyzed By: AH Prep Batch: 54106 QC Preparation: 2009-09-07 Prepared By: AH

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Cyanide	0.121	$_{ m mg/L}$	1	0.120	< 0.0110	101	62.6 - 132

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		Rec .		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Cyanide	0.125	mg/L	1	0.120	< 0.0110	104	62.6 - 132	3	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 209098

QC Batch: 63425Date Analyzed: 2009-09-10 Analyzed By: DS Prepared By: DS Prep Batch: 54137 QC Preparation: 2009-09-04

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec .	Limit
HMX	1.41	$\mu \mathrm{g/L}$	1	2.50	< 0.123	56	10 - 164
RDX	1.63	$\mu { m g}/{ m L}$	1	2.50	< 0.298	65	10 - 147
1,3,5-Trinitrobenzene	1.76	$\mu { m g}/{ m L}$	1	2.50	< 0.339	70	10 - 187
1,3-Dinitrobenzene	1.87	$\mu { m g}/{ m L}$	1	2.50	< 0.389	75	10 - 155
Nitrobenzene	1.87	$\mu { m g}/{ m L}$	1	2.50	< 0.379	75	10 - 156
Tetryl	1.80	$\mu { m g}/{ m L}$	1	2.50	< 0.413	72	10 - 158
TNT	1.93	$\mu { m g}/{ m L}$	1	2.50	< 0.464	77	21 - 114
4-Amino-DNT	2.19	$\mu { m g}/{ m L}$	1	2.50	< 0.319	88	80 - 120

 $continued \dots$

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matrix spikes continued . . .

		MS			Spike	Matrix		$\mathrm{Rec}.$
Param		Result	Units	Dil.	${f Amount}$	Result	$\mathrm{Rec}.$	Limit
2-Amino-DNT	13	1.96	$\mu \mathrm{g/L}$	1	2.50	< 0.391	78	80 - 120
2,6-DNT	14	1.64	$\mu { m g}/{ m L}$	1	2.50	< 0.323	66	80 - 120
2,4-DNT	15	1.96	$\mu { m g}/{ m L}$	1	2.50	< 0.366	78	80 - 120
2-NT		1.99	$\mu { m g}/{ m L}$	1	2.50	< 0.379	80	10 - 147
4-NT		1.81	$\mu { m g}/{ m L}$	1	2.50	< 0.398	72	10 - 161
3-NT		2.08	$\mu { m g}/{ m L}$	1	2.50	< 0.346	83	10 - 167

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

		MSD			Spike	Matrix		Rec .		RPD
Param		Result	Units	Dil.	${f Amount}$	Result	$\mathrm{Rec}.$	Limit	RPD	Limit
HMX		1.56	$\mu \mathrm{g/L}$	1	2.50	< 0.123	62	10 - 164	10	20
RDX		1.68	$\mu { m g/L}$	1	2.50	< 0.298	67	10 - 147	3	20
1,3,5-Trinitrobenzene		1.95	$\mu { m g/L}$	1	2.50	< 0.339	78	10 - 187	10	20
1,3-Dinitrobenzene		1.94	$\mu { m g/L}$	1	2.50	< 0.389	78	10 - 155	4	20
${ m Nitrobenzene}$		2.03	$\mu { m g/L}$	1	2.50	< 0.379	81	10 - 156	8	20
Tetryl		2.02	$\mu { m g/L}$	1	2.50	< 0.413	81	10 - 158	12	20
TNT		1.89	$\mu \mathrm{g/L}$	1	2.50	< 0.464	76	21 - 114	2	20
4-Amino-DNT		2.40	$\mu \mathrm{g/L}$	1	2.50	< 0.319	96	80 - 120	9	20
2-Amino-DNT		2.26	$\mu \mathrm{g/L}$	1	2.50	< 0.391	90	80 - 120	14	20
2,6-DNT	16	1.90	$\mu \mathrm{g/L}$	1	2.50	< 0.323	76	80 - 120	15	20
2,4-DNT		2.09	$\mu { m g/L}$	1	2.50	< 0.366	84	80 - 120	6	20
2-NT		1.97	$\mu \mathrm{g/L}$	1	2.50	< 0.379	79	10 - 147	1	20
4-NT		1.92	$\mu \mathrm{g/L}$	1	2.50	< 0.398	77	10 - 161	6	20
3-NT		1.92	$\mu { m g/L}$	1	2.50	< 0.346	77	10 - 167	8	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MS	MSD			Spike	MS	MSD	$\mathrm{Rec}.$
$\mathbf{Surrogate}$	Result	Result	Units	Dil.	${f Amount}$	$\mathrm{Rec.}$	$\mathrm{Rec}.$	Limit
1,2-Dinitrobenzene	3.49	3.68	$\mu { m g/L}$	1	2.5	140	147	10 - 222

Matrix Spike (MS-1) Spiked Sample: 209230

QC Batch: 63441 Date Analyzed: 2009-09-09 Analyzed By: AH Prep Batch: 54150 QC Preparation: 2009-09-09 Prepared By: AH

	MS			Spike	Matrix		Rec .
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Total Kjeldahl Nitrogen - N	44.5	${ m mg/L}$	1	50.0	2.52	84	61.2 - 118

¹³Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

¹⁴Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

 $^{^{15}}$ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

¹⁶ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

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	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Total Kjeldahl Nitrogen - N	46.2	mg/L	1	50.0	2.52	87	61.2 - 118	4	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208408

QC Batch: 63461 Date Analyzed: 2009-09-11 Analyzed By: RR
Prep Batch: 54153 QC Preparation: 2009-09-11 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit Dissolved Chromium 0.110 0.100 0.014 96 75 - 125 mg/L1

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSD Spike Matrix RPD Rec. Param Result Units Dil. Amount Result Rec. Limit RPD Limit Dissolved Chromium 20 0.112 mg/L0.100 0.014 75 - 125

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208427

QC Batch: 63485 Date Analyzed: 2009-09-11 Analyzed By: KV
Prep Batch: 54188 QC Preparation: 2009-09-11 Prepared By: KV

MSSpike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit Nitrate and Nitrite as N 19.2 mg/L200 0.200 14.52350 80 - 120

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

MSD RPD Spike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit RPD Limit Nitrate and Nitrite as N 18.7 200 0.200 9350 20 mg/L14.580 - 120

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208955

 $continued \dots$

¹⁷Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

¹⁸Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

${\it Work~Order:~9090211} \\ {\it HELSTF~Diesel~Spill~Groundwater}$

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matrix spikes continued								
•		MS			Spike	Matrix		$\mathrm{Rec}.$
Param		Result	Units	Dil .	${ m Amount}$	Result	$\mathrm{Rec.}$	Limit
		MS			Spike	Matrix		Rec.
Param		Result	Units	Dil.	$\overline{\mathrm{Amount}}$	Result	$\mathrm{Rec.}$	Limit
Chloride	19	129	$\mathrm{mg/L}$	5.56	139	6	88	90 - 110

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

		MSD			Spike	Matrix		Rec .		RPD
Param		Result	Units	Dil.	Amount	Result	Rec.	Limit	RPD	Limit
Chloride	20	128	mg/L	5.56	139	6	88	90 - 110	1	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208955

QC Batch: 63668 Date Analyzed: 2009-09-01 Analyzed By: JR
Prep Batch: 54356 QC Preparation: 2009-09-01 Prepared By: JR

MS Spike Matrix Rec. Param Result Units Dil. Amount Result Rec. Limit Sulfate 127mg/L5.56139 < 2.8090 - 110

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		$\mathrm{Rec}.$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Sulfate	127	mg/L	5.56	139	< 2.80	91	90 - 110	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Matrix Spike (MS-1) Spiked Sample: 208953

QC Batch: 63674 Date Analyzed: 2009-09-16 Analyzed By: JR
Prep Batch: 54363 QC Preparation: 2009-09-16 Prepared By: JR

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${ m Amount}$	Result	$\mathrm{Rec.}$	Limit
Bromide	10700	mg/L	2222	11100	<87.5	96	90 - 110

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Bromide	10800	$\mathrm{mg/L}$	2222	11100	<87.5	97	90 - 110	1	20

¹⁹Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

²⁰ Matrix spike recovery out of control limits due to matrix interference. Use LCS/LCSD to demonstrate analysis is under control.

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Matrix Spike (MS-1) Spiked Sample: 208953

QC Batch: 63674Prep Batch: 54363

Date Analyzed: 2009-09-16 QC Preparation: 2009-09-16 Analyzed By: JR Prepared By: JR

	MS			Spike	Matrix		$\mathrm{Rec}.$
Param	Result	Units	Dil.	${f Amount}$	Result	Rec .	Limit
Fluoride	10900	$\mathrm{mg/L}$	2222	11100	< 96.4	98	90 - 110

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

	MSD			Spike	Matrix		${ m Rec.}$		RPD
Param	Result	Units	Dil.	${ m Amount}$	Result	Rec.	Limit	RPD	Limit
Fluoride	10900	mg/L	2222	11100	<96.4	98	90 - 110	0	20

Percent recovery is based on the spike result. RPD is based on the spike and spike duplicate result.

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
Total Silver		mg/L	0.250	0.253	101	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	$\operatorname{Recovery}$	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${\bf Analyzed}$
Total Aluminum		$\mathrm{mg/L}$	1.00	1.01	101	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
Total Arsenic		$\mathrm{mg/L}$	2.00	2.02	101	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Work Order: 9090211 HELSTF Diesel Spill Groundwater

			CCVs True	CCVs Found	CCVs Percent	Percent Recovery	Date
Param	Flag	Units	Conc.	Conc.	Recovery	Limits	${\bf Analyzed}$
Total Barium		${ m mg/L}$	1.00	1.02	102	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

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			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Beryllium		mg/L	1.00	1.02	102	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Cadmium		$\mathrm{mg/L}$	1.00	1.04	104	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
Total Cobalt		$_{ m mg/L}$	1.00	0.992	99	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	Analyzed
Total Chromium		$\mathrm{mg/L}$	1.00	1.04	104	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Work Order: 9090211 HELSTF Diesel Spill Groundwater

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Copper		$_{ m mg/L}$	1.00	1.05	105	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

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			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Iron		$_{ m mg/L}$	1.00	1.04	104	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Manganese		$\mathrm{mg/L}$	1.00	1.01	101	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Molybdenum		$_{ m mg/L}$	1.00	0.992	99	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Nickel		mg/L	1.00	0.999	100	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Work Order: 9090211 HELSTF Diesel Spill Groundwater

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Phosphorous		m mg/L	5.00	4.93	99	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

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			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Lead		$\mathrm{mg/L}$	2.00	2.06	103	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	Analyzed
Total Antimony		$\mathrm{mg/L}$	2.00	2.04	102	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Selenium		$_{ m mg/L}$	1.00	1.02	102	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Thallium		$\mathrm{mg/L}$	5.00	5.10	102	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Work Order: 9090211 HELSTF Diesel Spill Groundwater

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Vanadium		mg/L	1.00	1.04	104	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

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			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	$\operatorname{Recovery}$	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
Total Zinc		$\mathrm{mg/L}$	1.00	1.08	108	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	$\operatorname{Recovery}$	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	Analyzed
Total Silver		$\mathrm{mg/L}$	0.125	0.129	103	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Aluminum		$_{ m mg/L}$	1.00	0.993	99	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Arsenic		$_{ m mg/L}$	1.00	0.981	98	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Work Order: 9090211 HELSTF Diesel Spill Groundwater

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STE Diogol Spill Croundwater	

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	Analyzed
Total Barium		m mg/L	1.00	1.04	104	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Beryllium		mg/L	1.00	1.00	100	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Cadmium		mg/L	1.00	1.01	101	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Cobalt		$\mathrm{mg/L}$	1.00	0.989	99	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Chromium		mg/L	1.00	1.02	102	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221 Date Analyzed: 2009-09-03 Analyzed By: RR

Work Order: 9090211 HELSTF Diesel Spill Groundwater

			0.0177	0.077	0.077	_	
			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	\mathbf{Found}	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Copper		mg/L	1.00	1.04	104	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

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			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Iron		m mg/L	1.00	1.02	102	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Manganese		$\mathrm{mg/L}$	1.00	0.996	100	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Molybdenum		$_{ m mg/L}$	1.00	0.985	98	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Nickel		mg/L	1.00	0.973	97	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Work Order: 9090211 HELSTF Diesel Spill Groundwater

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Phosphorous		mø/L	5.00	4 84	97	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

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			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Lead		m mg/L	1.00	0.966	97	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	Analyzed
Total Antimony		$\mathrm{mg/L}$	1.00	1.00	100	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Selenium		$_{ m mg/L}$	1.00	0.983	98	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Thallium		$\mathrm{mg/L}$	1.00	1.00	100	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Work Order: 9090211 HELSTF Diesel Spill Groundwater

			CCVs True	CCVs Found	CCVs Percent	Percent Recovery	Date
Param	Flag	Units	Conc.	Conc.	Recovery	Limits	${\bf Analyzed}$
Total Vanadium		$\mathrm{mg/L}$	1.00	1.04	104	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63221

Date Analyzed: 2009-09-03

Analyzed By: RR

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			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	$\operatorname{Recovery}$	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
Total Zinc		$_{ m mg/L}$	1.00	1.03	103	90 - 110	2009-09-03

Standard (ICV-1)

QC Batch: 63227

Date Analyzed: 2009-09-03

Analyzed By: TP

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc.	Conc.	Recovery	Limits	${ m Analyzed}$
Total Mercury		$\mathrm{mg/L}$	0.00100	0.000990	99	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63227

Date Analyzed: 2009-09-03

Analyzed By: TP

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Mercury		mg/L	0.00100	0.000960	96	90 - 110	2009-09-03

Standard (CCV-1)

QC Batch: 63239

Date Analyzed: 2009-09-03

Analyzed By: ER

Analyzed By: ER

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
GRO		$_{ m mg/L}$	1.00	1.07	107	80 - 120	2009-09-03

Standard (CCV-2)

QC Batch: 63239

Date Analyzed: 2009-09-03

 ${\it Work~Order:~9090211} \\ {\it HELSTF~Diesel~Spill~Groundwater}$

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			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	$\operatorname{Recovery}$	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
GRO		$_{ m mg/L}$	1.00	1.02	102	80 - 120	2009-09-03

Standard (CCV-1)

QC Batch: 63261 Date Analyzed: 2009-09-04 Analyzed By: MN

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
Phenol		m mg/L	60.0	55.1	92	80 - 120	2009-09-04
1,4-Dichlorobenzene (para)		m mg/L	60.0	59.4	99	80 - 120	2009-09-04
2-Nitrophenol		m mg/L	60.0	65.0	108	80 - 120	2009-09-04
2,4-Dichlorophenol		${ m mg/L}$	60.0	67.4	112	80 - 120	2009-09-04
${\it Hexachlorobutadiene}$		${ m mg/L}$	60.0	61.7	103	80 - 120	2009-09-04
4-Chloro-3-methylphenol	21	${ m mg/L}$	60.0	73.5	122	80 - 120	2009-09-04
2,4,6-Trichlorophenol		${ m mg/L}$	60.0	63.5	106	80 - 120	2009-09-04
${ m Acenaphthene}$		m mg/L	60.0	60.2	100	80 - 120	2009-09-04
Diphenylamine		m mg/L	60.0	60.4	101	80 - 120	2009-09-04
${ m Pentachlorophenol}$		m mg/L	60.0	58.4	97	80 - 120	2009-09-04
Fluoranthene		${ m mg/L}$	60.0	59.8	100	80 - 120	2009-09-04
Di-n-octylphthalate		m mg/L	60.0	63.6	106	80 - 120	2009-09-04
$\mathrm{Benzo}(\mathrm{a})\mathrm{pyrene}$		m mg/L	60.0	62.7	104	80 - 120	2009-09-04

					${ m Spike}$	$\operatorname{Percent}$	$\operatorname{Recovery}$
$\mathbf{Surrogate}$	Flag	Result	Units	$\operatorname{Dilution}$	${f Amount}$	Recovery	Limit
2-Fluorophenol		61.7	$\mathrm{mg/L}$	1	60.0	103	80 - 120
${ m Phenol-d5}$		56.5	${ m mg/L}$	1	60.0	94	80 - 120
${ m Nitrobenzene-d5}$		58.2	$\mathrm{mg/L}$	1	60.0	97	80 - 120
2-Fluorobiphenyl		58.2	$\mathrm{mg/L}$	1	60.0	97	80 - 120
2,4,6-Tribromophenol		70.7	${ m mg/L}$	1	60.0	118	80 - 120
Terphenyl-d14		57.7	${ m mg/L}$	1	60.0	96	80 - 120

Standard (CCV-1)

QC Batch: 63307 Date Analyzed: 2009-09-07 Analyzed By:

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
DRO		m mg/L	250	286	114	80 - 120	2009-09-07

Standard (CCV-2)

QC Batch: 63307 Date Analyzed: 2009-09-07 Analyzed By:

²¹Control analyte out of CCV control limits. Results biased high. •

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			H	ELSTF Diese	el Spill G	roundwater		
Param	Flag	$_{ m Units}$	CCVs True Conc.	CCV Foun Cond	d :.	CCVs Percent Recovery	Percent Recovery Limits	$egin{array}{c} { m Date} \ { m Analyzed} \end{array}$
DRO		$\mathrm{mg/L}$	250	262		105	80 - 120	2009-09-07
Standard	(CCV-1)							
QC Batch:	63327		Date	Analyzed:	2009-09-0	1	Analy	zed By: MD
Param		Flag	Units	${ m CCVs} \ { m True} \ { m Conc.}$	CCVs Found Conc.	$egin{array}{c} ext{CCVs} \ ext{Percent} \ ext{Recovery} \end{array}$	Percent Recovery Limits	$\begin{array}{c} {\rm Date} \\ {\rm Analyzed} \end{array}$
Hexavalent	Chromium		$\mathrm{mg/L}$	0.500	0.495	99	90 - 110	2009-09-01
Standard QC Batch:	(CCV-2) 63327		Date	Analyzed:	2009-09-0	1	${ m Analy}$	zed By: MD
				${ m CCVs} \ { m True}$	$\begin{array}{c} { m CCVs} \\ { m Found} \end{array}$	$rac{ ext{CCVs}}{ ext{Percent}}$	$egin{array}{c} ext{Percent} \ ext{Recovery} \end{array}$	Date
Param	CI.	Flag	Units	Conc.	Conc.	Recovery	Limits	Analyzed
Hexavalent	Chromium		mg/L	0.500	0.483	97	90 - 110	2009-09-01
Standard	(ICV-1)							
QC Batch:	63339		Date	Analyzed:	2009-09-0	01	Anal	yzed By: JG
Param	Flag	Units	$egin{array}{c} \mathrm{CCVs} \ \mathrm{True} \ \mathrm{Conc.} \end{array}$	CCV Foun Cond	d	CCVs Percent Recovery	Percent Recovery Limits	$egin{array}{c} { m Date} \ { m Analyzed} \end{array}$
pH	2 1006	s.u.	7.00	6.97		100	98 - 102	2009-09-01
Standard QC Batch:	(CCV-1) 63339		Date	$\label{eq:Analyzed} \textbf{Analyzed:}$	2009-09-0	01	Anal	yzed By: JG
			$\overset{ ext{CCVs}}{ ext{-}}$	CCV	s	$_{ m CCVs}$	Percent	_

Standard (CCV-2)

Flag

 Units

s.u.

 Param

 \overline{pH}

QC Batch: 63340 Date Analyzed: 2009-09-08 Analyzed By: KV

Found

 ${\rm Conc.}$

6.97

Percent

Recovery

100

Recovery

 ${\bf Limits}$

98 - 102

Date

 ${\bf Analyzed}$

2009-09-01

 ${\rm True}$

Conc.

7.00

Work Order: 9090211 HELSTF Diesel Spill Groundwater

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Organic Carbon		mø/L	50.0	49.8	100	80 - 120	2009-09-08

Standard (CCV-3)

QC Batch: 63340

Date Analyzed: 2009-09-08

Analyzed By: KV

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			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Organic Carbon		$\mathrm{mg/L}$	50.0	49.0	98	80 - 120	2009-09-08

Standard (ICV-1)

QC Batch: 63355

Date Analyzed: 2009-09-09

Analyzed By: RR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Calcium		m mg/L	50.0	51.5	103	90 - 110	2009-09-09

Standard (ICV-1)

QC Batch: 63355

Date Analyzed: 2009-09-09

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Potassium		$_{ m mg/L}$	50.0	50.8	102	90 - 110	2009-09-09

Standard (ICV-1)

QC Batch: 63355

Date Analyzed: 2009-09-09

Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Magnesium		mg/L	50.0	51.0	102	90 - 110	2009-09-09

Standard (ICV-1)

QC Batch: 63355

Date Analyzed: 2009-09-09

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HELSTF Diesel Spill Groundwater	

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Sodium		$\mathrm{mg/L}$	50.0	50.6	101	90 - 110	2009-09-09

Standard (CCV-1)

QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
Total Calcium		mg/L	50.0	51.1	102	90 - 110	2009-09-09

Standard (CCV-1)

QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	Analyzed
Total Potassium		$_{ m mg/L}$	50.0	51.1	102	90 - 110	2009-09-09

Standard (CCV-1)

QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	Analyzed
Total Magnesium		$_{ m mg/L}$	50.0	50.4	101	90 - 110	2009-09-09

Standard (CCV-1)

QC Batch: 63355 Date Analyzed: 2009-09-09 Analyzed By: RR

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Sodium		$_{ m mg/L}$	50.0	50.1	100	90 - 110	2009-09-09

Standard (ICV-1)

QC Batch: 63370 Date Analyzed: 2009-09-05 Analyzed By: AH

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Ammonia-N		$_{ m mg/L}$	5.00	4.93	99	85 - 115	2009-09-05

Standard (CCV-1)

QC Batch: 63370

Date Analyzed: 2009-09-05

Analyzed By: AH

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			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	$\operatorname{Recovery}$	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Ammonia-N		$_{ m mg/L}$	5.00	4.98	100	85 - 115	2009-09-05

Standard (ICV-1)

QC Batch: 63391

Date Analyzed: 2009-09-07

Analyzed By: AH

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Total Cyanide		m mg/L	0.120	0.118	98	85 - 115	2009-09-07

Standard (CCV-1)

QC Batch: 63391

Date Analyzed: 2009-09-07

Analyzed By: AH

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Cyanide		$_{ m mg/L}$	0.120	0.124	103	85 - 115	2009-09-07

Standard (ICV-1)

QC Batch: 63425

Date Analyzed: 2009-09-10

Analyzed By: DS

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	Analyzed
HMX		$\mu \mathrm{g/L}$	500	505	101	85 - 115	2009-09-10
RDX		$\mu { m g}/{ m L}$	500	478	96	85 - 115	2009-09-10
1,3,5-Trinitrobenzene		$\mu { m g}/{ m L}$	500	490	98	85 - 115	2009-09-10
1,3-Dinitrobenzene		$\mu { m g}/{ m L}$	500	507	101	85 - 115	2009-09-10
${ m Nitrobenzene}$		$\mu { m g}/{ m L}$	500	507	101	85 - 115	2009-09-10
Tetryl		$\mu { m g}/{ m L}$	500	485	97	85 - 115	2009-09-10
TNT		$\mu { m g}/{ m L}$	500	481	96	85 - 115	2009-09-10
4-Amino-DNT		$\mu { m g}/{ m L}$	500	518	104	85 - 115	2009-09-10
2-Amino-DNT		$\mu { m g}/{ m L}$	500	540	108	85 - 115	2009-09-10

 $continued \dots$

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TOTE DE LO ILO L	

standard continued							
			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
$\overline{2,6\text{-DNT}}$		$\mu \mathrm{g/L}$	500	466	93	85 - 115	2009-09-10
2,4-DNT		$\mu { m g}/{ m L}$	500	520	104	85 - 115	2009-09-10
2-NT		$\mu { m g}/{ m L}$	500	503	101	85 - 115	2009-09-10
4-NT		$\mu { m g}/{ m L}$	500	433	87	85 - 115	2009-09-10
3-NT		$\mu { m g}/{ m L}$	500	493	99	85 - 115	2009-09-10

					Spike	$\operatorname{Percent}$	Recovery
$\mathbf{Surrogate}$	Flag	Result	Units	Dilution	${f Amount}$	Recovery	Limit
1,2-Dinitrobenzene		456	$\mu \mathrm{g}/\mathrm{L}$	1	500	91	85 - 115

Standard (CCV-1)

 $QC\ Batch: \quad 63425$ Date Analyzed: 2009-09-10 Analyzed By: DS

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${ m Analyzed}$
HMX		$\mu \mathrm{g}/\mathrm{L}$	500	545	109	85 - 115	2009-09-10
RDX		$\mu { m g}/{ m L}$	500	461	92	85 - 115	2009-09-10
1,3,5-Trinitrobenzene		$\mu { m g}/{ m L}$	500	483	97	85 - 115	2009-09-10
1,3-Dinitrobenzene		$\mu { m g}/{ m L}$	500	516	103	85 - 115	2009-09-10
${ m Nitrobenzene}$		$\mu { m g}/{ m L}$	500	511	102	85 - 115	2009-09-10
Tetryl		$\mu { m g}/{ m L}$	500	467	93	85 - 115	2009-09-10
TNT		$\mu { m g}/{ m L}$	500	486	97	85 - 115	2009-09-10
4-Amino-DNT		$\mu { m g}/{ m L}$	500	538	108	85 - 115	2009-09-10
2-Amino-DNT		$\mu { m g}/{ m L}$	500	551	110	85 - 115	2009-09-10
2,6-DNT		$\mu { m g}/{ m L}$	500	502	100	85 - 115	2009-09-10
2,4-DNT		$\mu { m g}/{ m L}$	500	552	110	85 - 115	2009-09-10
2-NT		$\mu { m g}/{ m L}$	500	535	107	85 - 115	2009-09-10
4-NT		$\mu { m g}/{ m L}$	500	500	100	85 - 115	2009-09-10
3-NT		$\mu { m g}/{ m L}$	500	507	101	85 - 115	2009-09-10

					Spike	Percent	Recovery
$\operatorname{Surrogate}$	Flag	Result	Units	Dilution	${f Amount}$	Recovery	Limit
1,2-Dinitrobenzene		491	$\mu { m g/L}$	1	500	98	85 - 115

Standard (ICV-1)

QC Batch: 63441 Date Analyzed: 2009-09-09 Analyzed By: AH

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc.	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Total Kjeldahl Nitrogen - N		$_{ m mg/L}$	5.00	5.21	104	85 - 115	2009-09-09

Work Order: 9090211 HELSTF Diesel Spill Ground Report Date: September 18, 2009 Page Number: 77 of 80

		H	HELSTF Die	sel Spill Grou	ındwater		
Standard (CCV-1)							
QC Batch: 63441		Date	e Analyzed:	2009-09-09		Analy	zed By: AH
Param	\mathbf{Fl}	ag Unit	CCVs True s Conc.	$\begin{array}{c} { m CCVs} \\ { m Found} \\ { m Conc.} \end{array}$	CCVs Percent Recovery	Percent Recovery Limits	$egin{array}{c} { m Date} \ { m Analyzed} \end{array}$
Total Kjeldahl Nitrogen		$\frac{\rm mg/l}{\rm mg}$		4.93	99	85 - 115	2009-09-09
		- 07					
Standard (ICV-1)							
QC Batch: 63461		Date	e Analyzed:	2009-09-11		Analy	zed By: RR
			$ ext{CCVs}$	CCVs	$_{ m CCVs}$	Percent	ъ.
Danara	Elo.	Units	${ m True} \ { m Conc.}$	Found	Percent	Recovery	$egin{array}{c} { m Date} \\ { m Analyzed} \end{array}$
Param Dissolved Chromium	Flag	mg/L	1.00	Conc. 1.03	Recovery 103	Limits 90 - 110	2009-09-11
Dissolved Chromitum		mg/L	1.00	1.03	103	90 - 110	2009-09-11
Standard (CCV-1)							
QC Batch: 63461		Date	e Analyzed:	2009-09-11		Analy	zed By: RR
Param	Flag	Units	$egin{array}{c} \mathrm{CCVs} \\ \mathrm{True} \\ \mathrm{Conc.} \end{array}$	$\begin{array}{c} {\rm CCVs} \\ {\rm Found} \\ {\rm Conc.} \end{array}$	CCVs Percent Recovery	Percent Recovery Limits	$\begin{array}{c} \text{Date} \\ \text{Analyzed} \end{array}$
Dissolved Chromium		mg/L	1.00	1.02	102	90 - 110	2009-09-11
Standard (ICV-1) QC Batch: 63473		Date	e Analyzed:	2009-09-03		m Analy z	zed By: MD
			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	${ m Units}$	$\operatorname{Conc.}$	Conc.	Recovery	Limits	Analyzed
Total Dissolved Solids		$\mathrm{mg/L}$	1000	994	99	90 - 110	2009-09-03
Standard (CCV-1) QC Batch: 63473			e Analyzed:	2009-09-03		Analyz	zed By: MD
			$ ext{CCVs}$	CCVs	$_{ m CCVs}$	Percent	ъ
Param	Flag	Units	${ m True} \ { m Conc.}$	Found Conc.	$egin{array}{l} ext{Percent} \ ext{Recovery} \end{array}$	$egin{array}{c} ext{Recovery} \ ext{Limits} \end{array}$	$egin{array}{c} { m Date} \\ { m Analyzed} \end{array}$
1 611 61111	riag	Omto	COHe.	COHe.	Trecovery	типпо	Anaryzeu

Standard (ICV-1)

Total Dissolved Solids

QC Batch: 63485 Analyzed By: KV Date Analyzed: 2009-09-11

989

99

90 - 110

2009-09-03

1000

mg/L

-							
			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${\bf Analyzed}$
Nitrate and Nitrite as N		mg/L	0.200	0.188	94	85 - 115	2009-09-11

Standard (CCV-1)

QC Batch: 63485

Date Analyzed: 2009-09-11

Analyzed By: KV

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			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Nitrate and Nitrite as N		$_{ m mg/L}$	0.200	0.201	100	85 - 115	2009-09-11

Standard (ICV-1)

QC Batch: 63527

Date Analyzed: 2009-09-10

Analyzed By: JG

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc.	Conc .	Recovery	Limits	Analyzed
Hydroxide Alkalinity		mg/L as $CaCo3$	0.00	<1.00		90 - 110	2009-09-10
Carbonate Alkalinity		mg/L as $CaCo3$	0.00	240		90 - 110	2009-09-10
Bicarbonate Alkalinity		mg/L as CaCo3	0.00	10.0		90 - 110	2009-09-10
Total Alkalinity		mg/L as $CaCo3$	250	250	100	90 - 110	2009-09-10

Standard (CCV-1)

QC Batch: 63527

Date Analyzed: 2009-09-10

Analyzed By: JG

			CCVs	CCVs	CCVs	Percent	
			True	Found	Percent	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Hydroxide Alkalinity		mg/L as CaCo3	0.00	<1.00		90 - 110	2009-09-10
Carbonate Alkalinity		mg/L as $CaCo3$	0.00	240		90 - 110	2009-09-10
Bicarbonate Alkalinity		mg/L as $CaCo3$	0.00	10.0		90 - 110	2009-09-10
Total Alkalinity		mg/L as $CaCo3$	250	250	100	90 - 110	2009-09-10

Standard (CCV-1)

QC Batch: 63668

Date Analyzed: 2009-09-01

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Chloride		m mg/L	25.0	22.9	92	90 - 110	2009-09-01

HELSTF Diesel Spill Groundwater

Work Order: 9090211 Page Number: 79 of 80

Analyzed By: JR

Standard (C	CV-1)
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QC Batch: 63668 Date Analyzed: 2009-09-01

> CCVsCCVs CCVs Percent

True Found Percent Recovery Date Flag Param Units Conc. Conc. Recovery Limits Analyzed 22.6 2009-09-01 Sulfate mg/L25.0 90 - 110 90

Standard (CCV-2)

QC Batch: 63668 Date Analyzed: 2009-09-01 Analyzed By: JR

CCVsCCVsCCVsPercent True Found Percent Recovery Date Param Flag Units Conc. Conc. Recovery Limits Analyzed 90 - 110 Chloride 23.1 2009-09-01 mg/L25.092

Standard (CCV-2)

QC Batch: 63668 Date Analyzed: 2009-09-01 Analyzed By: JR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	$\operatorname{Conc.}$	Recovery	Limits	${ m Analyzed}$
Sulfate		$_{ m mg/L}$	25.0	22.7	91	90 - 110	2009-09-01

Standard (CCV-1)

QC Batch: 63674 Date Analyzed: 2009-09-16 Analyzed By: JR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	$\operatorname{Conc.}$	Conc .	Recovery	Limits	${\bf Analyzed}$
Bromide		m mg/L	5.00	4.90	98	90 - 110	2009-09-16

Standard (CCV-1)

QC Batch: 63674 Date Analyzed: 2009-09-16 Analyzed By: JR

			CCVs	CCVs	CCVs	$\operatorname{Percent}$	
			True	Found	$\operatorname{Percent}$	Recovery	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	Analyzed
Fluoride		m mg/L	5.00	5.06	101	90 - 110	2009-09-16

Standard (CCV-2)

 $QC\ Batch:\ 63674$ Date Analyzed: 2009-09-16 Analyzed By: JR

Work Order: 9090211 HELSTF Diesel Spill Groundwater

			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	$\operatorname{Recovery}$	Date
Param	Flag	Units	$\operatorname{Conc.}$	$\operatorname{Conc.}$	Recovery	Limits	${\bf Analyzed}$
Bromide		$_{ m mg/L}$	5.00	5.08	102	90 - 110	2009-09-16

Standard (CCV-2)

QC Batch: 63674

Date Analyzed: 2009-09-16

Analyzed By: JR

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			CCVs	CCVs	CCVs	Percent	
			True	Found	$\operatorname{Percent}$	$\operatorname{Recovery}$	Date
Param	Flag	Units	Conc .	Conc .	Recovery	Limits	${ m Analyzed}$
Fluoride		$_{ m mg/L}$	5.00	5.16	103	90 - 110	2009-09-16

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